

Exhibit F



Handbook of Fine Chemicals and Laboratory Equipment

Aldrich

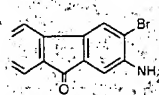
TP
202
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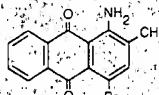
2003-2004

US
\$

W 361.15 mp 282°(dec.)	1g	22.80
50050 IRRITANT	5g	75.80
sulfonic acid, sodium		
266		
mp 148-149°	1g	33.70
50050 IRRITANT		
174.12 mp 213-214°	250mg	34.30
T	1g	93.90
174.12 mp 243-244°	100mg	49.30
81C IRRITANT		
5] FW 316.16	5g	26.00
15(12), 1693D IRRITANT	25g	86.50
	100g	210.80
81-50-5]	25g	22.20
	100g	58.40
FW 187.04 mp 73-76°	1g	21.50
	5g	90.40
W 187.04 mp 88-95°	1g	19.20
	5g	75.20
FW 187.04 mp 92-96°	1g	24.40
	5g	83.30
35-3] (5-bromo-6-methyl-2,828A Safety 2,145D	10g	29.20
no-1-naphthylamine)	5g	44.10
(1), 1252A Safety 2,146A	25g	144.10
1] H ₂ NC ₆ H ₂ (Br)(NO ₂)CN	1g	10.30
	5g	33.70
218.02 mp 205-208°	1g	21.60
VT	5g	69.90
pyridine-3-carbonitrile	1g	29.80
FW 255.13 mp 183-187°	5g	30.10
	25g	133.10
mp >350° Beil. 26(5), 16,103	1g	55.10
	5g	182.50
mp 136-138° Beil. 22,431	5g	21.60
	25g	72.10
2' mp 88-91° IRRITANT	5g	51.20
	25g	170.70



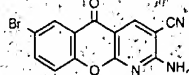
29,979-0



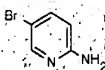
24,669-7



37,688-4

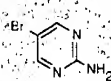


40,243-5

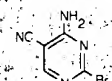


12,285-8

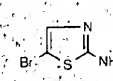
30,352-6	2-Amino-5-bromopyrimidine, 98% [7752-82-1] FW 174.01 mp 241-243°	1g	28.00
	Beil. 25(4), 2125 Safety 2,146C R&S 1(2), 2575D IRRITANT	5g	92.50
	Intermediate for the preparation of sulfanilamides and amino acids containing the pyrimidine ring system. The products are potential antiviral agents. J. Chem. Ed. 1985, 62, 905.		
56,384-6	4-Amino-2-bromopyrimidine-5-carbonitrile, 95% [94741-70-5] FW 199.01	1g	44.50
	mp 254°(dec.) IRRITANT		
12,802-3	2-Amino-5-bromothiazole monohydrobromide, 97% [3034-22-8] FW 259.96	5g	38.20
	mp 165°(dec.) FT-NMR 1(3), 115A FT-IR 1(2), 645D R&S 1(2), 2387J	25g	124.00
30,708-4	(R)-(-)-2-Amino-1-butanol, 98% [5856-63-3] C ₂ H ₅ CH(NH ₂)CH ₂ OH FW 89.14 mp -2°	1g	10.90
	bp 172-174° n _D ²⁰ 1.4520 d 0.947 Fp 180°F(82°C) [α] _D ²⁰ -10° (neat) Beil. 4,291 Merck	5g	35.20
	Index 13,438 FT-NMR 1(1), 538B FT-IR 1(1), 337C Safety 2,147A R&S 1(1), 363D		
	CORROSIVE		
	96% ee/GLC		
A4,380-4	2-Amino-1-butanol, 97% [96-20-8] C ₂ H ₅ CH(NH ₂)CH ₂ OH FW 89.14 mp -2°	5mL	16.20
	bp 176-178° n _D ²⁰ 1.4510 d 0.943 Fp 184°F(84°C) [α] _D ²⁰ 0° (c=2, C ₂ H ₅ OH) Beil. 4,291	100mL	24.20
	Merck. Index 13,426 FT-NMR 1(1), 539A FT-IR 1(1), 337C Safety 2,147B R&S 1(1), 363F	500mL	77.90
	RTECS# EK9625000 CORROSIVE		
13,252-7	(S)-(+)-2-Amino-1-butanol, 98+% [5856-62-2] C ₂ H ₅ CH(NH ₂)CH ₂ OH FW 89.14	1g	22.70
	bp 172-174° n _D ²⁰ 1.4520 d 0.944 Fp 175°F(79°C) [α] _D ²⁰ +10° (neat) Beil. 4,291 Merck	5g	75.10
	Index 13,438 FT-NMR 1(1), 538C FT-IR 1(1), 337D Safety 2,146D R&S 1(1), 363E		
	CORROSIVE		
	96% ee/GLC		
17,833-0	4-Amino-1-butanol, 98% [13325-10-5] H ₂ N(CH ₂) ₄ OH FW 89.14 bp 206° n _D ²⁰ 1.4610	1g	23.80
	d 0.967 Fp 226°F(107°C) Beil. 4,291 FT-NMR 1(1), 538A FT-IR 1(1), 338D	5g	78.80
	Safety 2,147D R&S 1(1), 363C RTECS# EK9700000 CORROSIVE HYGROSCOPIC		
07191	4-Amino-1-butanol, purum, ~97% (GC) [13325-10-5]	1mL	23.50
		5mL	80.00
		25mL	335.80
29,460-8	3-Amino-2-butenethioamide, 98% [62069-87-8] CH ₃ C(NH ₂)=CHCSNH ₂ FW 116.19	25g	52.20
	mp 135-138° Beil. 4(4), 2846 FT-NMR 1(1), 1335C Safety 2,148A R&S 1(1), 971G		
	IRRITANT		
36,084-8	(Z)-2-Amino-N-[(1-ethoxyethoxy)carbonyl]-1-methylethoxyimino]-4-thiazoleacetic acid, 98% [86299-47-0] FW 329.38 mp 180°(dec.) FT-NMR 1(3), 496C	5g	23.40
	R&S 1(2), 26911 MOISTURE-SENSITIVE IRRITANT	25g	76.40
38,747-9	3-Amino-N-butylbenzenesulfonamide, tech., 90% [143173-93-7]	5g	27.50
	H ₂ NC ₆ H ₄ SO ₂ NH(CH ₂) ₃ CH ₃ FW 228.32 mp 73-77° R&S 1(2), 2237N IRRITANT	25g	84.70
	6-[N-(4-Aminobutyl)-N-ethylamino]-2,3-dihydro-1,4-phthalazinedione, see 26,906-9; N-(4-Aminobutyl)-N-ethylsoluminol, page 73		
26,906-9	N-(4-Aminobutyl)-N-ethylsoluminol [66612-29-1] [6-[N-(4-aminobutyl)-N-ethylamino]-2,3-dihydro-1,4-phthalazinedione] FW 276.34 mp 259-260°	25mg	64.50
	FT-NMR 1(2), 1416B Safety 2,148B R&S 1(2), 2041H LIGHT-SENSITIVE	100mg	213.80
	Efficient chemiluminescent NH ₂ -coupling reagent for detection of a wide variety of proteins down to the picomole range. Nature 1979, 279, 646. Methods Enzymol. 1978, 57, 424.		
	May contain up to 15% water.		
	(4-Aminobutyl)guanidine sulfate, see 10,144-3; Agmatine sulfate, page 43		
19,328-3	2-Amino-4-tert-butylphenol, 98% [1199-46-8] (CH ₃) ₃ CC ₆ H ₃ (NH ₂)OH FW 165.24	10g	36.00
	mp 160-163° FT-NMR 1(2), 515C FT-IR 1(1), 1227B Safety 2,148C R&S 1(1), 1417A	50g	88.10
	IRRITANT		
52,440-9	2-Amino-5-tert-butyl-1,3,4-thiadiazole, 97% [39222-73-6] FW 157.23 mp 183-187°	5g	46.40
	IRRITANT		



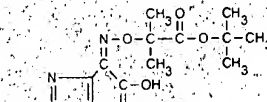
30,352-6



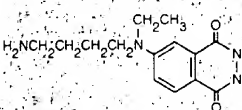
56,384-6



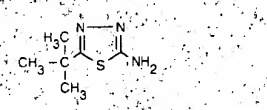
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36,084-8

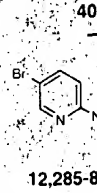
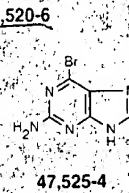
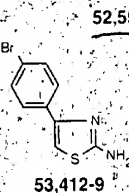
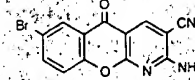
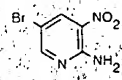
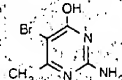
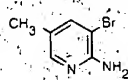
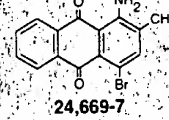
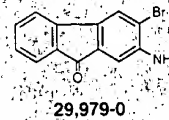
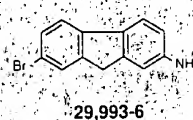
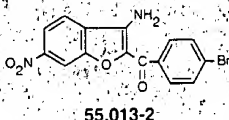


26,906-9



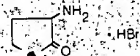
52,440-9

55,013-2	3-Amino-2-(4-bromobenzoyl)-6-nitrobenzofuran, 97% FW 361.15 mp 282°(dec.)	1g	US \$ 22.80
NEW	IRRITANT	5g	75.80
29,993-6	1-Amino-4-bromo-9,10-dihydro-9,10-dioxo-2-anthracenesulfonicacid, sodium salt, see 45,606-3. Bromaminic acid, sodium salt page 266	1g	33.70
Beil. 12(2).780 Safety 2.144C R&S 1(1).1449J RTECS# LL5150050	IRRITANT		
29,979-0	2-Amino-3-bromo-9-fluorenone, 95% [52086-09-6] FW 274.12 mp 213-214°	250mg	34.30
FT-NMR 1(2).907B Safety 2.144D R&S 1(2).1681B	IRRITANT	1g	93.90
30,015-2	2-Amino-7-bromo-9-fluorenone, 98% [58557-63-4] FW 274.12 mp 243-244°	100mg	49.30
Beil. 14(2).69 FT-NMR 1(2).907C Safety 2.145A R&S 1(2).1681C	IRRITANT		
24,669-7	1-Amino-4-bromo-2-methylantraquinone, 99% [81-50-5] FW 316.16	5g	26.00
★ mp 245°(dec.) Beil. 14.223 FT-IR 1(2).91B Safety 2.145C R&S 1(2).1693D	IRRITANT	25g	86.50
		100g	210.80
24,670-0	1-Amino-4-bromo-2-methylantraquinone, tech., 95% [81-50-5]	25g	22.20
★		100g	58.40
54,307-1	2-Amino-3-bromo-5-methylpyridine, 98% [17282-00-7] FW 187.04 mp 73-76°	1g	21.50
NEW	IRRITANT	5g	90.40
52,553-7	2-Amino-5-bromo-3-methylpyridine, 97% [3430-21-5] FW 187.04 mp 88-95°	1g	19.20
NEW	IRRITANT	5g	75.20
54,840-5	6-Amino-3-bromo-2-methylpyridine, 97% [42753-71-9] FW 187.04 mp 92-96°	1g	24.40
NEW ★	IRRITANT	5g	83.30
20,520-6	2-Amino-5-bromo-6-methyl-4-pyrimidinol, 99% [6307-35-3] (5-bromo-6-methyl-isocytosine) FW 204.03 mp 244-246° Beil. 24.350 FT-IR 1(2).828A Safety 2.145D R&S 1(2).2579D	10g	29.20
	IRRITANT		
A4,340-5	1-Amino-4-bromonaphthalene, 97% [2298-07-9] (4-bromo-1-naphthylamine) BrC ₁₀ H ₆ NH ₂ FW 222.09 mp 102-103° Beil. 12.1257 FT-IR 1(1).1252A Safety 2.146A R&S 1(1).1445M	5g	44.10
	IRRITANT	25g	144.10
51,692-9	2-Amino-3-bromo-5-nitrobenzonitrile, 97% [17601-94-4] H ₂ NC ₆ H ₂ (Br)(NO ₂)CN FW 242.03 mp 180-185°	1g	10.30
NEW ★	IRRITANT	5g	33.70
37,688-4	2-Amino-5-bromo-3-nitropyridine, 97% [6945-68-2] FW 218.02 mp 205-208° Beil. 22(1).631 FT-NMR 1(3).306B R&S 1(2).2533B	1g	21.60
	IRRITANT	5g	69.90
40,243-5	2-Amino-7-bromo-5-oxo-5H-[1]benzopyrano[2,3-b]pyridine-3-carbonitrile, 99% [206658-80-2] FW 316.12 mp >300°	1g	29.80
	IRRITANT		
53,412-9	2-Amino-4-(4-bromophenyl)thiazole, 97% [2103-94-8] FW 255.13 mp 183-187°	5g	30.10
NEW	IRRITANT	25g	133.10
47,525-4	2-Amino-6-bromopurine, 95% [82499-03-4] FW 214.03 mp >350° Beil. 26(5).16.103	1g	55.10
		5g	182.50
12,285-8	2-Amino-5-bromopyridine, 97% [1072-97-5] FW 173.02 mp 136-138° Beil. 22.431 FT-IR 1(2).772B Safety 2.146B R&S 1(2).2529D	5g	21.60
	IRRITANT	25g	72.10
52,174-4	2-Amino-6-bromopyridine, 98% [19798-81-3] FW 173.02 mp 88-91°	5g	51.20
	IRRITANT	25g	170.70

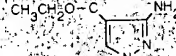


Aminobutyr

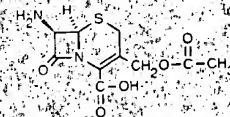
			US \$
A4,415-0	4-Aminobutylaldehyde diethyl acetal, tech. 90% [6346-09-4] (4,4-diethoxybutylamine) $H_2N(CH_2)_3CH(OC_2H_5)_2$ FW 161.25 bp 196° n _D 1.4270 d 0.933 Fp 145°F (62°C) Beil. 4.319 FT-NMR 1(1), 533B FT-IR 1(1), 331D Safety 2.149A R&S 1(1), 359E IRRITANT	1g 5g 25g 100g	11.30 36.70 121.60 329.60
11,612-2	(R)-(-)-2-Aminobutyric acid, 99% [2623-91-8] (D-2-aminobutyric acid) $C_2H_5CH(NH_2)CO_2H$ FW 103.12 mp >300° [α] _D -7.94° (c=4, H ₂ O) Beil. 4.408 FT-NMR 1(1), 871C FT-IR 1(1), 573A R&S 1(1), 655C 99% ee/GLC	1g 5g	27.10 116.50
16,266-3	DL-2-Aminobutyric acid, 99% [2835-81-6] (DL-2-aminobutyric acid) $C_2H_5CH(NH_2)CO_2H$ FW 103.12 mp 291° (dec.) Beil. 4.408 Merck Index 13.439 FT-NMR 1(1), 871B FT-IR 1(1), 572C R&S 1(1), 655B	25g 100g	19.80 54.80
23,438-9	(S)-(+)-2-Aminobutyric acid, 98% [1492-24-6] (L-2-aminobutyric acid) $C_2H_5CH(NH_2)CO_2H$ FW 103.12 [α] _D +20.4° (c=2, 5N HCl) Beil. 4.408 Merck Index 13.439 FT-NMR 1(1), 872A FT-IR 1(1), 572D R&S 1(1), 655D 99% ee/GLC	1g 5g	25.30 83.90
A4,420-7	DL-3-Aminobutyric acid, 97% [2835-82-7] (DL-3-aminobutyric acid) $CH_3CH(NH_2)CH_2CO_2H$ FW 103.12 mp 189° (dec.) Beil. 4.412 Merck Index 13.440 FT-NMR 1(1), 877B FT-IR 1(1), 577B R&S 1(1), 657O	5g 25g	26.00 99.70
A4,440-1	4-Aminobutyric acid, 97% [56-12-2] (γ-aminobutyric acid, GABA) $H_2N(CH_2)_3CO_2H$ FW 103.12 mp 195° (dec.) Beil. 4.413 Merck Index 13.429 FT-IR 1(1), 577C R&S 1(1), 659A RTECS# ES6300000 IRRITANT Important inhibitory neurotransmitter! Reacts with isothiocyanates to produce thioureas which have antifungal activity? (1) Neurochem. Res. 1993; 18: 365; Chem. Abstr. 1994, 120: 225805w (2) Eur. J. Med. Chem. 1991; 26: 723	25g 100g 500g	11.20 26.60 105.30
A4,450-9	DL-Amino-γ-butyrolactone hydrobromide, 99% [6305-38-0] FW 182.02 mp 221° (dec.) [α] _D 25° (c=20, H ₂ O) Beil. 18.601 FT-NMR 1(1), 1139C FT-IR 1(1), 700C R&S 1(1), 807O	1g 10g 25g	13.90 76.60 152.50
47,142-9	(S)-(-)-Amino-γ-butyrolactone hydrobromide, 99% [15295-77-9] FW 182.02 mp 225° (dec.) [α] _D -21° (c=1, H ₂ O) IRRITANT	1g 5g	22.00 72.40
46,247-0	(R)-(+)-Amino-γ-butyrolactone hydrochloride, 97% [104347-13-9] FW 137.57 mp 220-224° [α] _D +28° (c=1, H ₂ O) IRRITANT	1g 5g	49.50 166.60
45,922-4	(S)-(-)-Amino-γ-butyrolactone hydrochloride, 97% [2185-03-7] FW 137.57 mp 215-220° (dec.) [α] _D -27.8° (c=1, H ₂ O) Beil. 18(5), 11.314 IRRITANT	50mg 250mg 1g	14.40 54.10 106.70
A4,460-6	6-Aminocaproic acid, 98% [60-32-2] (6-aminohexanoic acid) $H_2N(CH_2)_5CO_2H$ FW 131.18 mp 210° (dec.) Beil. 4.434 Merck Index 13.431 FT-IR 1(1), 578D Safety 2.149B R&S 1(1), 659G RTECS# MO6300000 IRRITANT Hemostat! Used as a lysine analogue to characterize binding sites in plasminogen. 2,3 (1) Thromb. Res. 1994; 73: 419; Chem. Abstr. 1994, 120: 208282x (2) Biochemistry 1994; 33: 3599 (3) ibid. 1993; 32: 13681	5g 100g 500g 1kg	6.40 17.80 58.70 108.10
21,770-0	DL-2-Aminocaprylic acid, 99% [644-90-6] (DL-2-aminooctanoic acid) $CH_3(CH_2)_5CH(NH_2)CO_2H$ FW 159.23 mp 260° (dec.) Beil. 4.461 FT-NMR 1(1), 876C FT-IR 1(1), 577A R&S 1(1), 657M	5g 25g	43.30 144.00
85,529-4	8-Aminocaprylic acid, 99% [1002-57-9] (8-aminooctanoic acid) $H_2N(CH_2)_7CO_2H$ FW 159.23 mp 194° (dec.) FT-NMR 1(1), 879B FT-IR 1(1), 579A R&S 1(1), 659I	1g 5g	34.70 139.10
A4,500-9	3-Amino-4-carbethoxypyrazole, 98% [6994-25-8] (ethyl 3-amino-4-pyrazole carboxylate) FW 155.16 mp 105-107° FT-IR 1(2), 610A R&S 1(2), 2347N CORROSIVE	1g 5g	15.50 51.70
43,951-7	1-(Aminocarbonyl)-1-cyclopropanecarboxylic acid, 97% [6914-74-5] $H_2NCO_2CH_2CO_2H$ FW 129.12 mp 185° (dec.) Beil. 9(2), 512 IRRITANT	1g 5g	33.50 110.30
19,114-0	7-Aminocephalosporanic acid, 99% [957-68-6] (7-ACA) FW 272.28 mp >300° [α] _D +90° (c=0.5, KH ₂ PO ₄ /trace NaOH) Merck Index 13.433 FT-NMR 1(1), 1317B FT-IR 1(1), 809B R&S 1(1), 949F Potent inhibitor of bacterial (S. aureus) β-lactamase. J. Med. Chem. 1974; 17: 342	1g 5g	34.90 123.20
51,610-4	3-Amino-4-chloroacetanilide, 97% [51867-83-5] $CH_3CONHC_6H_4(NH_2)Cl$ FW 184.62 mp 172-176° IRRITANT	25g	140.30
54,824-3	4-Amino-5-chloro-o-anisic acid; see 34.087 1,4-Amino-5-chloro-2-methoxy benzoic acid page 76	5g 25g	31.30 104.10



A4,450-9



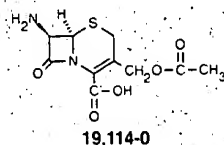
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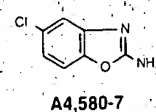
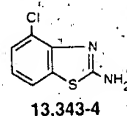
19,114-0

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f) (4,4-diethoxy- ... 70 d 0.933 safety 2,149A	1g 5g 25g 100g	11.30 36.70 121.60 329.60
rylic acid) Beil. 4,408	1g 5g	27.10 116.50
cid) Index 13,439	25g 100g	19.80 54.80
rylic acid) 108 Merck	1g 5g	25.30 83.90
cid) ck Index 13,440	5g 25g	26.00 99.70
ABA) lex 13,429	25g 100g 500g	11.20 26.60 105.30
to produce thioureas Chem. Abstr. 1994,		
W 182.02 9C FT-IR 1(1),700C	1g 10g 25g	13.90 76.60 152.50
-77-9] FW 182.02 5g	1g 5g	22.00 72.40
47-13-9] FW 137.57 ... 5g	1g 5g	49.50 166.60
03-7] FW 137.57 TANT	50mg 250mg 1g	14.40 54.10 106.70
d) H ₂ N(CH ₂) ₅ CO ₂ H ... 1(1),578D	5g 100g 500g	6.40 17.80 58.70
in plasminogen ^{2,3} (2) Biochemistry.	1kg	108.10
noic acid) FT-NMR 1(1),876C	5g 25g	43.30 144.00
acid) H ₂ N(CH ₂) ₅ CO ₂ H &S 1(1),659I	1g 5g	34.70 139.10
mino-4-pyrazole- 2),2347N CORROSIVE	1g 5g	15.50 51.70
914-74-5] TANT	1g 5g	33.50 110.30
N 272.28 mp >300° NMR 1(1),1317B	1g 5g	34.90 123.20
n. 1974, 17, 342. tC ₆ H ₃ (NH ₂)Cl	25g	140.30
chloro-2-methoxy-		
CONH ₂ FW 170.60 ... 25g	5g 25g	31.30 104.10

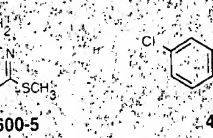
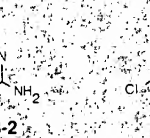
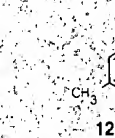
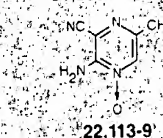
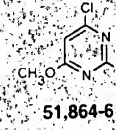
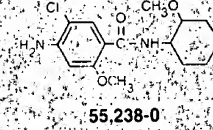
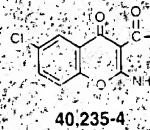
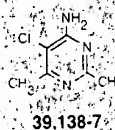
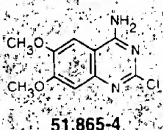


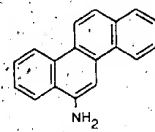
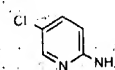
30,856-0	4-Amino-6-chloro-1,3-benzenedisulfonamide, 98% [121-30-2] ★ H ₂ NC ₆ H ₂ (Cl)(SO ₂ NH ₂) ₂ FW 285.73 mp 257-261° Beil. 14(4),2810 Merck Index 13,2086 FT-NMR 1(2),1645A R&S 1(2),2251A MOISTURE-SENSITIVE IRRITANT	5g 100g	20.60 54.90
14,297-2	3-Amino-4-chlorobenzenesulfonyl fluoride, 98% [368-72-9] H ₂ NC ₆ H ₃ (Cl)SO ₂ F FW 209.63 mp 62-64° FT-IR 1(2),524A Safety 2,150D R&S 1(2),2233E CORROSIVE MOISTURE-SENSITIVE	1g	26.50
54,185-0	2-Amino-4-chlorobenzenethiol, 96% [1004-00-8] NH ₂ C ₆ H ₃ (Cl)SH FW 159.64 (NEW) mp 56-61° CORROSIVE STENCH	5g 25g	39.10 130.20
36,227-1	2-Amino-3-chlorobenzoic acid, 95% [6388-47-2] (3-chloroanthranilic acid) H ₂ NC ₆ H ₃ (Cl)CO ₂ H FW 171.58 mp 189-191° Beil. 14(4),1071 FT-NMR 1(2),1098B R&S 1(2),1799A IRRITANT	1g	28.40
A4,546-7	2-Amino-4-chlorobenzoic acid, 98% [89-77-0] (4-chloroanthranilic acid) H ₂ NC ₆ H ₃ (Cl)CO ₂ H FW 171.58 mp 231-233° Beil. 14,365 FT-NMR 1(2),1119B FT-IR 1(2),218A R&S 1(2),1807I RTECS# DG1570000 IRRITANT	25g 100g	30.90 81.50
A4,547-5	2-Amino-5-chlorobenzoic acid, 98% [635-21-2] (5-chloroanthranilic acid) ★ H ₂ NC ₆ H ₃ (Cl)CO ₂ H FW 171.58 mp 210-212° Beil. 14,365 FT-NMR 1(2),1120B FT-IR 1(2),218D R&S 1(2),1807L IRRITANT	1g 10g	24.50 83.00
37,804-6	2-Amino-5-chlorobenzoic acid, tech., 90% [635-21-2] (5-chloroanthranilic acid) ★ H ₂ NC ₆ H ₃ (Cl)CO ₂ H	5g 25g	37.60 124.50
42,262-2	2-Amino-6-chlorobenzoic acid, 98% [2148-56-3] (6-chloroanthranilic acid) H ₂ NC ₆ H ₃ (Cl)CO ₂ H FW 171.58 mp 158-160° Beil. 14(3),966 IRRITANT	1g 10g	15.80 81.50
42,599-0	3-Amino-2-chlorobenzoic acid, 97% [108679-71-6] H ₂ NC ₆ H ₃ (Cl)CO ₂ H FW 171.58 mp 154-160° IRRITANT	100mg 500mg	21.70 70.60
23,936-4	3-Amino-4-chlorobenzoic acid, 98% [2840-28-0] H ₂ NC ₆ H ₃ (Cl)CO ₂ H FW 171.58 ... mp 214-215° Beil. 14,412 FT-NMR 1(2),1119A FT-IR 1(2),217D Safety 2,151A R&S 1(2),1807H IRRITANT	25g 100g	33.90 121.40
21,771-9	4-Amino-2-chlorobenzoic acid, 97% [2457-76-3] H ₂ NC ₆ H ₃ (Cl)CO ₂ H FW 171.58 ... ★ mp 211°(dec.) Beil. 14,438 FT-NMR 1(2),1119C FT-IR 1(2),218B Safety 2,151B R&S 1(2),1807J RTECS# DG1575000 IRRITANT	1g 5g	14.60 50.00
53,249-5	5-Amino-2-chlorobenzoic acid, 97% [89-54-3] H ₂ NC ₆ H ₃ (Cl)CO ₂ H FW 171.58 (NEW) ★ mp 184-188° IRRITANT	25g	38.30
24,825-8	5-Amino-2-chlorobenzoic acid, tech., 85% [89-54-3] H ₂ NC ₆ H ₃ (Cl)CO ₂ H ★	5g 100g 500g	14.40 46.70 118.40
40,528-0	2-Amino-4-chlorobenzonitrile, 99% [38487-86-4] (4-chloroanthranilonitrile) ★ H ₂ NC ₆ H ₃ (Cl)CN FW 152.58 mp 161-162° Beil. 14(3),954 IRRITANT	10g 50g	28.50 93.90
17,433-5	2-Amino-5-chlorobenzonitrile, 98% [5922-60-1] (5-chloroanthranilonitrile) ★ H ₂ NC ₆ H ₃ (Cl)CN FW 152.58 mp 96-99° bp 132-135°/0.5mm Fp >230°F(110°C) FT-IR 1(2),458A Safety 2,151D R&S 1(2),2137C IRRITANT	10g	26.00
27,901-3	4-Amino-2-chlorobenzonitrile, 99% [20925-27-3] H ₂ NC ₆ H ₃ (Cl)CN FW 152.58 mp 117-119° R&S 1(2),2137D IRRITANT	50g	87.60
A4,556-4	2-Amino-5-chlorobenzophenone, 98% [719-59-5] H ₂ NC ₆ H ₃ (Cl)COC ₆ H ₅ FW 231.68 mp 98-100° Beil. 14,79 FT-IR 1(2),70D Safety 2,152A R&S 1(2),1675H RTECS# PC4933500 IRRITANT	25g 100g 500g	11.90 28.20 83.90
13,343-4	2-Amino-4-chlorobenzothiazole, 97% [19952-47-7] FW 184.65 mp 203-205° ★ FT-NMR 1(3),207B FT-IR 1(2),703C R&S 1(2),2457B RTECS# DL1225000 IRRITANT	5g 25g 100g	42.10 131.70 370.50
13,608-5	2-Amino-6-chlorobenzothiazole, 99% [95-24-9] FW 184.65 mp 199-201° ★ Beil. 27(2),230 FT-NMR 1(3),207C R&S 1(2),2457C RTECS# DL1575000 TOXIC IRRITANT	5g 25g	22.70 89.60
	Aminochlorobenzotrifluoride, see Chloro(trifluoromethyl)aniline		
A4,580-7	2-Amino-5-chlorobenzoxazole, 97% [61-80-3] (zoxazolamine) FW 168.58 mp 181-184° Merck Index 13,10249 FT-NMR 1(3),199B FT-IR 1(2),698D Safety 2,153A R&S 1(2),2451F RTECS# DM4550000 IRRITANT	1g 5g 25g	18.30 60.80 240.30
21,861-8	2-(3-Amino-4-chlorobenzoyl)benzoic acid, 95+% [118-04-7] ★ H ₂ NC ₆ H ₃ (Cl)COC ₆ H ₄ CO ₂ H FW 275.69 mp 171-173° Beil. 14,661 FT-NMR 1(2),1167B FT-IR 1(2),240C R&S 1(2),1829G IRRITANT	50g	34.50
55,015-9	2-Amino-2-(4-chlorobenzoyl)-6-nitrobenzofuran, 97% FW 316.70 mp 262°(dec.) ... (NEW) IRRITANT	1g 5g	20.90 69.40



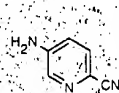
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33,953-9	2-Amino-5-chlorobenzyl alcohol, 99% [37585-25-4] $\text{H}_2\text{NC}_6\text{H}_3(\text{Cl})\text{CH}_2\text{OH}$ FW 157.60 mp 107-109° FT-NMR 1(2), 517C R&S 1(1), 1417H IRRITANT	1g 5g	19.60 66.30
51,865-4	4-Amino-2-chloro-6,7-dimethoxyquinazoline, 95% [23680-84-4] FW 239.66 mp 260°(dec.) Beil. 25(5), 13,350 IRRITANT	25g	213.60
39,138-7	4-Amino-5-chloro-2,6-dimethylpyrimidine, 96% [2858-20-0] FW 157.60 mp 164-166° Beil. 24,90 IRRITANT	5g 25g	73.60 244.60
39,721-0	2-Amino-4-chlorodiphenylamine, see 32,752-2, N-(4-Chlorophenyl)-1,2-phenylenediamine page 464		
39,721-0	2-Amino-5-chloro-2'-fluorobenzophenone, 98% [784-38-3] $\text{H}_2\text{NC}_6\text{H}_3(\text{Cl})\text{COC}_6\text{H}_4\text{F}$ FW 249.67 mp 95-98° IRRITANT Employed in the syntheses of benzotriazepines ¹ and diazepam-related benzodiazepines ² (1) J. Chem. Res. Synop. 1991; 2. (2) J. Am. Chem. Soc. 1990, 112, 3969	5g 25g	22.00 73.00
40,235-4	2-Amino-6-chloro-3-formylchromone, 97% [68301-77-9] (2-amino-6-chloro-4-oxo-4H-1-benzopyran-3-carboxaldehyde) FW 223.62 mp >300° IRRITANT	1g	26.30
34,087-1	4-Amino-5-chloro-2-methoxybenzoic acid, 97% [7206-70-4] (4-amino-5-chloro-o-anisic acid) $\text{H}_2\text{NC}_6\text{H}_3(\text{Cl})(\text{OCH}_3)\text{CO}_2\text{H}$ FW 201.61 mp 206°(dec.) FT-NMR 1(2), 1150A R&S 1(2), 1821K IRRITANT	5g 25g	13.80 45.80
55,238-0	(4-Amino-5-chloro-2-methoxy)-N-(3-methoxy(4-piperidyl))benzamide, 97% [84946-16-7] FW 313.78 mp 190-193° IRRITANT	25g	160.50
51,864-6	2-Amino-4-chloro-6-methoxypyrimidine, 95% [5734-64-5] FW 159.58 mp 168-171° Beil. 25,7 IRRITANT	10g 50g	51.20 120.10
	4-Amino-2'-chloro-5'-methylbenzanilide, see 27,405-4, Fast Red Violet LB base page 894		
22,113-9	3-Amino-6-(chloromethyl)-2-pyrazinecarbonitrile 4-oxide, 97% [40127-89-7] FW 184.59 mp 144-146° FT-IR 1(2), 917D Safety 2, 154A R&S 1(2), 2683D IRRITANT	1g 5g	29.80 98.80
12,288-2	2-Amino-4-chloro-6-methylpyrimidine, 97% [5600-21-5] FW 143.58 mp 183-186° Beil. 24,84 FT-IR 1(2), 823C R&S 1(2), 2575I RTECS# UV6329600	5g 25g	29.20 96.50
A4,600-5	4-Amino-6-chloro-2-(methylthio)pyrimidine, 97% [1005-38-5] FW 175.64 mp 130-132° Beil. 25,11 FT-NMR 1(3), 382C FT-IR 1(2), 826A R&S 1(2), 2577I	1g 5g 25g	29.40 93.40 347.40
A4,640-4	1-Amino-4-chloronaphthalene, 98% [4684-12-2] (4-chloro-1-naphthylamine) $\text{ClC}_{10}\text{H}_6\text{NH}_2$ FW 177.63 mp 98-100° Beil. 12,1255 FT-IR 1(1), 1251D Safety 2, 154B R&S 1(1), 1445L IRRITANT	1g 5g	22.00 72.30
24,641-7	2-Amino-4-chloro-5-nitrophenol, 98% [6358-07-2] $\text{H}_2\text{NC}_6\text{H}_3(\text{Cl})(\text{NO}_2)\text{OH}$ FW 188.57 mp 225° Beil. 13,392 FT-NMR 1(2), 773A FT-IR 1(1), 1380D Safety 2, 155A R&S 1(1), 1601A RTECS# SJ5736000 IRRITANT	50g 250g	34.00 152.50
53,038-7	2-Amino-4-chloro-6-nitrophenol, 97% [6358-08-3] $\text{ClC}_6\text{H}_3(\text{NH}_2)(\text{NO}_2)\text{OH}$ FW 188.57 mp 158-162° IRRITANT	5g 25g	45.50 151.50
	2-Amino-4-chloro-5-nitrotoluene, see 20,120-0, 5-Chloro-4-nitro-o-toluidine page 457		
	2-Amino-6-chloro-4-oxo-4H-1-benzopyran-3-carboxaldehyde, see 40,235-4 2-Amino-6-chloro-3-formylchromone page 76		
40,231-1	2-Amino-7-chloro-5-oxo-5H-[1]benzopyrano[2,3-b]pyridine-3-carbonitrile, 97% [68302-10-3] FW 271.67 mp >300° IRRITANT	1g	28.10
C4,440-0	2-Amino-4-chlorophenol, 97% [95-85-2] (5-chloro-2-hydroxyaniline) $\text{H}_2\text{NC}_6\text{H}_3(\text{Cl})\text{OH}$ FW 143.57 mp 139-143° Beil. 13,383 FT-NMR 1(2), 517B FT-IR 1(1), 1228A Safety 2, 789A R&S 1(1), 1417F RTECS# SJ5700000 IRRITANT	5g 100g 500g	13.00 27.30 90.70
55,222-4	2-Amino-5-chlorophenol, 97% [28443-50-7] $\text{H}_2\text{NC}_6\text{H}_3(\text{Cl})\text{OH}$ FW 143.57 mp 145-153° IRRITANT	5g 25g	22.80 75.80
55,248-8	3-Amino-4-chlorophenol, 97% [16026-77-0] $\text{H}_2\text{NC}_6\text{H}_3(\text{Cl})\text{OH}$ FW 143.57 mp 151-154° IRRITANT	1g 5g	28.30 94.20

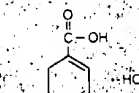


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40.231-1

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A4,710-9	4-Aminocinnamic acid hydrochloride, 97%, predominantly <i>trans</i> [54057-95-3]	1g	14.40
★	H ₂ NC ₆ H ₄ CH=CHCO ₂ H·HCl FW 199.64 mp 265-270°(dec.) Beil. 14.521	5g	47.20
	FT-NMR 1(2),1052A FT-IR 1(2),181C R&S 1(2),1775N IRRITANT	25g	148.70
33,657-2	2-Amino- <i>m</i> -cresol, 96% [2835-97-4] (2-amino-3-methylphenol) H ₂ NC ₆ H ₃ (CH ₃)OH	5g	44.20
	FW 123.16 mp 149-152° Beil. 13.589 FT-NMR 1(2),494B R&S 1(1),1405B IRRITANT	25g	116.50
14,490-8	2-Amino- <i>p</i> -cresol, 97% [95-84-1] (2-amino-4-methylphenol) H ₂ NC ₆ H ₃ (CH ₃)OH	50g	17.40
★	FW 123.16 mp 135-137° Beil. 13.601 FT-NMR 1(2),515A FT-IR 1(1),1227A	250g	57.20
	Safety 2,157A R&S 1(1),1415L RTECS# SJ6078000 IRRITANT		
22,506-1	3-Amino- <i>o</i> -cresol, 95% [53222-92-7] (3-amino-2-methylphenol)	250mg	29.40
	H ₂ NC ₆ H ₃ (CH ₃)OH FW 123.16 mp 129-130° Beil. 13.579 FT-NMR 1(2),514B	1g	85.20
	FT-IR 1(1),1226C Safety 2,157B R&S 1(1),1415J IRRITANT		
14,489-4	4-Amino- <i>m</i> -cresol, 97% [2835-99-6] (4-amino-3-methylphenol) H ₂ NC ₆ H ₃ (CH ₃)OH	5g	34.40
★	FW 123.16 mp 176-179° Beil. 13.593 FT-NMR 1(2),516B FT-IR 1(1),1227D	25g	114.00
	Safety 2,157D R&S 1(1),1417C RTECS# GO6826000 IRRITANT		
22,508-8	5-Amino- <i>o</i> -cresol, 97% [2835-95-2] (5-amino-2-methylphenol) H ₂ NC ₆ H ₃ (CH ₃)OH	5g	24.80
★	FW 123.16 mp 160-162° Beil. 13.574 RTECS# SJ6090000 IRRITANT	25g	81.90
14,491-6	6-Amino- <i>m</i> -cresol, 98% [2835-98-5] (2-amino-5-methylphenol) H ₂ NC ₆ H ₃ (CH ₃)OH	10g	26.00
	FW 123.16 mp 159-162° Beil. 13.590 FT-NMR 1(2),516A FT-IR 1(1),1227C	50g	90.80
	Safety 2,158A R&S 1(1),1417B IRRITANT		
11,764-1	3-Aminocrotonitrile, 96%, mixture of <i>cis</i> and <i>trans</i> [1118-61-2] (diacetonitrile)	5g	13.50
★	CH ₃ C(NH ₂)=CHCN FW 82.11 Beil. 3.660 FT-NMR 1(1),1383C FT-IR 1(1),861A	100g	29.30
	Safety 2,158B R&S 1(1),1009A LACHRYMATOR	500g	104.30
	5-Amino-6-cyano-2-pyrazinylmethyl acetate, see 86,162-6, 5-Acetoxyethyl-2-amino-3-cyanopyrazine page 17		
	3-Amino-4-cyanopyrazole, see 15,304-4, 3-Amino-4-pyrazolecarbonitrile page 108		
55,192-9	2-Amino-5-cyanopyridine, 97% [4214-73-7] FW 119.13 mp 159-163° IRRITANT	25g	65.50
NEW		100g	181.80
53,890-6	5-Amino-2-cyanopyridine, 96% [55338-73-3] FW 119.13 mp 148-152° IRRITANT	25g	119.40
NEW		100g	331.70
	(<i>R</i>)- <i>trans</i> -Amino-1,4-cyclohexadiene-1-acetic acid, see 47,138-0, (<i>R</i>)-(-)-2-(2,5-Dihydrophenyl)glycine page 678		
35,963-7	5-Amino-1,3-cyclohexadiene-1-carboxylic acid hydrochloride, 97%	10mg	36.60
	[59556-17-1] FW 175.62 mp 203°(dec.) FT-NMR 1(1),885A R&S 1(1),663F	50mg	121.40
	RTECS# GU4775000		
21,869-3	1-Amino-1-cyclohexanecarboxylic acid, 98% [2756-85-6] H ₂ NC ₆ H ₁₀ CO ₂ H	1g	12.90
★	FW 143.19 mp >300° Beil. 14.299 FT-NMR 1(1),884C FT-IR 1(1),584D R&S 1(1),663E	5g	28.80
	RTECS# GU8393000		
40,485-3	<i>cis</i> -4-Amino-1-cyclohexanecarboxylic acid, 95% [3685-23-2] H ₂ NC ₆ H ₁₀ CO ₂ H	500mg	26.60
	FW 143.19 mp 283°(dec.) Beil. 14.301 IRRITANT	1g	44.30
43,805-7	(<i>S</i>)-(+)-Aminocyclohexanepropionic acid hydrate, tech. [307310-72-1]	1g	25.00
	(3-cyclohexyl-L-alanine) C ₆ H ₁₁ CH ₂ CH(NH ₂)CO ₂ H·xH ₂ O FW 171.24 mp 234-237°	5g	87.10
	[α] _D ²⁵ +12.0° (c=1; 1N NaOH) IRRITANT		
22,257-7	<i>trans</i> -2-Aminocyclohexanol hydrochloride, 99% [5456-63-3] H ₂ NC ₆ H ₁₀ OH·HCl	5g	58.10
	FW 151.64 mp 172-175° Beil. 13(3),708 FT-NMR 1(1),542A FT-IR 1(1),315D	25g	190.50
	Safety 2,159A R&S 1(1),365F HYGROSCOPIC		
26,376-1	<i>trans</i> -4-Aminocyclohexanol hydrochloride, 97% [50910-54-8] H ₂ NC ₆ H ₁₀ OH·HCl	25g	58.20
	FW 151.64 mp 225-227° Beil. 13(2),159 FT-IR 1(1),316A Safety 2,159B R&S 1(1),365G	100g	159.50
	IRRITANT HYGROSCOPIC		
23,937-2	3-Amino-2-cyclohexen-1-one, 99% [5220-49-5] NH ₂ C ₆ H ₇ O FW 111.15	5g	24.60
	mp 131-133° FT-IR 1(1),460C Safety 2,159C R&S 1(1),483E IRRITANT	25g	82.00
52,232-5	2-Amino- <i>N</i> -cyclohexyl- <i>N</i> -methylbenzenesulfonamide, 97% [70693-59-3]	25g	58.10
★	H ₂ NC ₆ H ₄ SO ₂ N(C ₆ H ₁₁)CH ₃ FW 268.38 mp 98-101°		
46,343-4	2-Amino- <i>N</i> -cyclohexyl- <i>N</i> -methylbenzylamine, 98% [57365-08-9]	5g	17.10
	H ₂ NC ₆ H ₄ CH ₂ N(C ₆ H ₁₁)CH ₃ FW 218.34 mp 44-47° IRRITANT	25g	55.90
43,226-1	(<i>S</i>)-(+)-2-Amino-3-cyclohexyl-1-propanol hydrochloride, 97% [117160-99-3]	1g	21.30
	C ₆ H ₁₁ CH ₂ CH(NH ₂)CH ₂ OH·HCl FW 193.72 mp 230°(dec.) [α] _D ²⁵ +2.6° (c=1, CH ₃ OH)	5g	71.50
	IRRITANT		



53,890-6

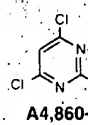
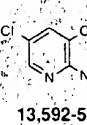
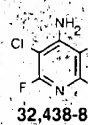
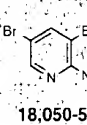
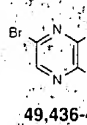
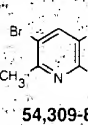
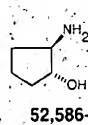
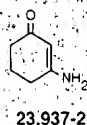


35,963-7

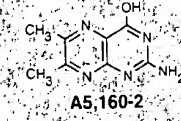
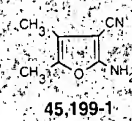
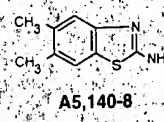
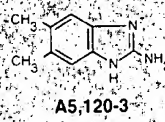
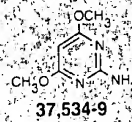
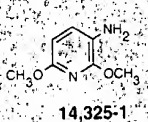
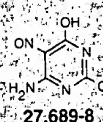


23,937-2

		US\$			US\$
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405B IRRITANT	1g 5g 25g	14.40 47.20 148.70	A4,810-5	1-Amino-1-cyclopentanecarboxylic acid, 97% [52-52-8] (cycloleucine)	5g 72.10 10g 119.60
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405B IRRITANT	5g 25g	44.20 116.50	19,227-9	1-Amino-1-cyclopentanemethanol, 97% [10316-79-7] (cycloleucinol)	1g 51.40
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405A	50g 250g	17.40 57.20	52,586-3	trans-2-Aminocyclopentanol hydrochloride, 97% [31775-67-4] FW 101.45 mp 191-196° [3160° (meat)]	5g 118.40
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405B	250mg 1g	29.40 85.20	30,408-5	1-Amino-1-cyclopropanecarboxylic acid hydrochloride, 97% [68781-13-5] 11-NC ₃ H ₆ (CO ₂ H)Cl FW 137.57 mp 231° (dec.) FT-NMR 1(1), 384A Safety 2, 160B R&S 1(1), 663B, IRRITANT	500mg 164.50
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405D	5g 25g	34.40 114.00	46,346-9	2-Amino-6H-dibenzol[b,d]pyran-6-one, see 30,023-3, 6-Amino-3,4-benzocoumarin page 69	
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405E	5g 25g	24.80 81.90	46,346-9	2-Amino-3,5-dibromobenzaldehyde, 97% [50910-55-9] H ₂ NC ₆ H ₃ (Br) ₂ CHO	5g 42.00 25g 139.20
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405F	10g 50g	26.00 90.80	46,346-9	4-Amino-3,5-dibromobenzenesulfonic acid, sodium salt, see 26,814-3, 3,5-Dibromosulfanilic acid, sodium salt page 591	
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405G	5g 25g	13.50 29.30 104.30	54,309-8	2-Amino-3,5-dibromo-6-methylpyridine, 97% [91872-10-5] FW 265.93 mp 143.5-148.5° IRRITANT	1g 41.70 5g 138.80
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405H	100g 500g	29.30 104.30	32,392-6	4-Amino-2,6-dibromophenol, 96% [609-21-2] H ₂ NC ₆ H ₃ (Br) ₂ OH FW 266.93 mp 195-196° Boil 13.517 FT-NMR 1(2), 532B R&S 1(1), 1425I IRRITANT	5g 17.00 25g 61.40
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405I	5g 25g	13.50 29.30 104.30	49,436-4	2-Amino-3,5-dibromopyrazine, 98% [24241-18-7] FW 252.91 mp 114-117° Boil 25(5)10, 146 IRRITANT	1g 54.50
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405J	5g 25g	13.50 29.30 104.30	18,050-5	2-Amino-3,5-dibromopyridine, 97% [35486-42-1] FW 251.92 mp 104-105° Boil 22.431 FT-IR 1(2), 773D Safety 2, 161A IRRITANT	5g 31.30
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405K	5g 25g	13.50 29.30 104.30	2-Amino-3,5-dichlorobenzoic acid, see 05,600-4, 3,5-Dichloroanthranilic acid page 604		
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405L	5g 25g	13.50 29.30 104.30	54,598-8	4-Amino-3,5-dichlorobenzoic acid, 98% [56961-25-2] H ₂ NC ₆ H ₃ (Cl) ₂ CO ₂ H	5g 22.10 25g 103.60
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405M	5g 25g	13.50 29.30 104.30	10,515-5	2-Amino-2,5-dichlorobenzophenone, 99% [2958-36-3] H ₂ NC ₆ H ₃ (Cl) ₂ COC ₆ H ₅	10g 38.50 50g 101.10
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405N	5g 25g	13.50 29.30 104.30	32,438-8	4-Amino-3,5-dichloro-2,6-difluoropyridine, 99% [2840-00-8] FW 198.99 mp 112-114° FT-NMR 1(3), 302B R&S 1(2), 2529M IRRITANT	1g 11.90 5g 39.20
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405O	5g 25g	13.50 29.30 104.30	38,579-4	6-Amino-2,4-dichloro-3-methylphenol hydrochloride, tech., 90% [39549-31-0] H ₂ NC ₆ H ₃ (Cl) ₂ (CH ₃)OH·HCl FW 228.51 R&S 1(1), 1427F IRRITANT	5g 15.80 25g 53.00
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405P	5g 25g	13.50 29.30 104.30	A4,840-7	4-Amino-2,6-dichlorophenol, 98% [5930-28-9] H ₂ NC ₆ H ₃ (Cl) ₂ OH FW 178.02 mp 167-170° Boil 13.512 FT-NMR 1(2), 528B FT-IR 1(1), 1231B Safety 2, 161D R&S 1(1), 1423G IRRITANT	25g 41.10 100g 121.70
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405Q	5g 25g	13.50 29.30 104.30	1-(4-Amino-3,5-dichlorophenyl)-ethanone, see 53,614-8, 3,5-Dichloro-4-aminoac etophenone page 603		
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405R	5g 25g	13.50 29.30 104.30	13,592-5	2-Amino-3,5-dichloropyridine, 97% [4214-74-8] FW 163.01 mp 81-83° Boil 22.430 FT-NMR 1(3), 301B FT-IR 1(2), 773C Safety 2, 162A R&S 1(2), 2529I IRRITANT	25g 88.70 100g 245.30
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405S	5g 25g	13.50 29.30 104.30	A4,860-1	2-Amino-4,6-dichloropyrimidine, 99% [56-05-3] FW 164.00 mp 219-222° Boil 24.81 FT-IR 1(2), 823B Safety 2, 162B R&S 1(2), 2575G RTECS# UV6260485 IRRITANT	10g 63.40 100g 369.80
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405T	5g 25g	13.50 29.30 104.30	21,773-5	5-Amino-4,6-dichloropyrimidine, 97% [5413-85-4] FW 164.00 mp 145-148° Boil 25(4), 2139 FT-IR 1(2), 826B Safety 2, 162C R&S 1(2), 2577J MOISTURE-SENSITIVE IRRITANT	1g 37.90 5g 148.00 25g 590.50
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405U	5g 25g	13.50 29.30 104.30	14,226-3	4-Amino-N,N-diethylamino- <i>o</i> -cresol dihydrochloride, 97% [6297-14-9] H ₂ NC ₆ H ₃ (CH ₂ N(C ₂ H ₅) ₂) ₂ OH·2HCl FW 267.20 mp 215-218° (dec.) FT-NMR 1(2), 628A FT-IR 1(1), 1297C R&S 1(1), 1499J	5g 43.10 25g 139.00
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405V	5g 25g	13.50 29.30 104.30	4-Amino-N-(2-diethylaminoethyl)benzamide hydrochloride, see 22,296-8, Procainamide hydrochloride page 1584		
11-NC ₆ H ₄ (CH ₂) ₃ OH 1405W	5g 25g	13.50 29.30 104.30	A4,880-6	2-Amino-5-diethylaminopentane, 97% [140-80-7] (N ¹ , N ¹ -diethyl-1,4-pentane diamine) (C ₂ H ₅) ₂ N(CH ₂) ₃ CH(NH ₂)CH ₃ FW 158.29 nD 1.4420 d 0.817 Fp 155°F (68°C) Merck Index 13.6756 FT-NMR 1(1), 498B FT-IR 1(1), 308A Safety 2, 162D R&S 1(1), 329J RTECS# XE1900000 CORROSIVE	25g 24.60 100g 55.00



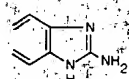
3-Amino-7-(diethylamino)-5-phenylphenazinium chloride, see 30.750-5		
Methylene Violet 3RAX page 1241		
3-Amino-N,N-diethyl-4-methoxybenzenesulfonamide, see 20.128-6. Fast Red ITR page 894		
2-Amino-4,5-difluorobenzoic acid, see 39.976-0. 4,5-Difluoroanthranilic acid page 662		
Amino-2,3-dihydro-1,4-phthalazinedione; see Aminophthalhydrazide		
27,689-8	6-Amino-2,4-dihydroxy-5-nitrosopyrimidine, 97% [5442-24-0] FW 156.10	1g 65.30 5g 270.30
Beil. 24.509 Safety 2.163B R&S 1(2).2585B IRRITANT		
A5,040-1	2-Amino-4,6-dihydroxypyrimidine, 98% [56-09-7] (2-amino-4,6-pyrimidinediol) FW 127.10 mp >300° Beil. 24.468 FT-IR 1(2).828D R&S 1(2).2579G	25g 34.40 100g 94.90
4-Amino-2,6-dihydroxypyrimidine, see A5,060-6. 6-Aminouracil page 113		
2-Amino-3,5-diiodobenzoic acid, see 28.348-7. 3,5-Diiodoanthranilic acid page 688		
A5,070-3	4-Amino-3,5-diiodobenzoic acid, 85% [2122-61-4] H ₂ NC ₆ H ₂ (I) ₂ CO ₂ H FW 388.93 mp >300° Beil. 14.439 FT-NMR 1(2).1149B FT-IR 1(2).232D R&S 1(2).1821I RTECS# DG2310000 IRRITANT	25g 60.80
4'-Amino-2',5'-dimethoxybenzanilide, see 20.154-5. Fast Blue RR page 893		
25,204-2	2-Amino-4,5-dimethoxybenzoic acid, 98% [5653-40-7] (6-aminoveratric acid; 4,5-dimethoxyanthranilic acid) H ₂ NC ₆ H ₂ (OCH ₃) ₂ CO ₂ H FW 197.19 mp 173°(dec.) Beil. 14.635 FT-NMR 1(2).1150B FT-IR 1(2).233A Safety 2.163C R&S 1(2).1821L IRRITANT	5g 32.30
36,111-9	2-Amino-4,5-dimethoxybenzonitrile, 96% [26961-27-3] H ₂ NC ₆ H ₂ (OCH ₃) ₂ CN FW 178.19 mp 98-99° Beil. 14(3).1624 FT-NMR 1(2).1541A R&S 1(2).2141H IRRITANT	5g 34.70
14,325-1	3-Amino-2,6-dimethoxypyridine monohydrochloride, 95% [80789-72-6] FW 190.63 mp 212° FT-IR 1(2).774A Safety 2.163D R&S 1(2).2529N IRRITANT	1g 61.70 5g 198.10
37,534-9	2-Amino-4,6-dimethoxypyrimidine, 98% [36315-01-2] FW 155.16 mp 94-96° Beil. 25(4).3485 FT-NMR 1(3).381B R&S 1(2).2575K IRRITANT	5g 18.70 25g 61.70
37,535-7	4-Amino-2,6-dimethoxypyrimidine, 97% [3289-50-7] FW 155.16 mp 149-152° R&S 1(2).2575L IRRITANT	5g 35.60
4'-Amino-2',3'-dimethylazobenzene, see 12.156-8. Fast Garnet GBC base page 893		
A5,120-3	2-Amino-5,6-dimethylbenzimidazole, 97% [29096-75-1] FW 161.21 mp 220-223°(dec.) FT-NMR 1(3).184C FT-IR 1(2).692B	5g 38.50 25g 141.00
A5,140-8	2-Amino-5,6-dimethylbenzothiazole, 97% [29927-08-0] FW 178.26 mp 185-189° FT-IR 1(2).703B R&S 1(2).2455N IRRITANT	5g 53.70
2-Amino-3,3-dimethylbutanoic acid, see tert-Leucine		
2-Amino-3,3-dimethyl-1-butanol, see 40.773-9. (S)-tert-Leucinol page 1123		
36,832-6	3-Amino-5,5-dimethyl-2-cyclohexen-1-one, 98% [873-95-0] H ₂ NC ₆ H ₅ (CH ₃) ₂ (=O) FW 139.20 mp 167-170° Beil. 7.560 FT-NMR 1(1).696C R&S 1(1).483F IRRITANT	5g 32.70
55,791-9	(S)-(+)-2-Amino-3,3-dimethyl-1,1-diphenylbutane, 99% (CH ₃) ₂ CCH(NH ₂)CH(C ₆ H ₅) ₂ FW 253.39 mp 50-53° Fp >230°F (110°C) [α] _D ²⁰ +24° (c=1, CHCl ₃) A Product of Onyx Scientific, U.K.	500mg 96.60
55,790-0	(S)-(-)-2-Amino-3,3-dimethyl-1,1-diphenyl-1-butanol, 99% [144054-70-6] (CH ₃) ₂ CCH(NH ₂)C(C ₆ H ₅) ₂ OH FW 269.39 mp 134-137° [α] _D ²⁰ -171° (c=1, CHCl ₃) MOISTURE-SENSITIVE IRRITANT A Product of Onyx Scientific, U.K.	1g 103.60 5g 344.70
45,199-1	2-Amino-4,5-dimethyl-3-furancarboxitrile, 95% [5117-88-4] FW 136.15 mp 163-169° Beil. 18(5).12.200	1g 27.30 10g 151.40
A5,160-2	2-Amino-6,7-dimethyl-4-hydroxypteridine, 97% [611-55-2] (2-amino-6,7-dimethyl-4-pteridinol) FW 191.19 mp >300° FT-IR 1(2).895C	500mg 72.80 1g 121.30



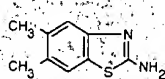
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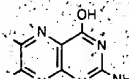
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anthranilic acid		
hydrazide		
-O) FW 156.10	1g	65.30
	5g	270.30
4,6-pyrimidinediol	25g	34.40
9G	100g	94.90
racil page 113		
anthranilic acid		
z(l)CO ₂ H FW 388.93	25g	60.80
1(2),1821L		
blue RR page 893		
aminoveratric acid,	5g	32.30
97.19 mp 173°(dec.)		
R&S 1(2),1821L		
IC ₆ H ₂ (OCH ₃) ₂ CN	5g	34.70
1(2),2141H IRRITANT		
[80789-72-6]	1g	61.70
29N IRRITANT	5g	198.10
55.16 mp 94-96°	5g	18.70
	25g	61.70
5.16 mp 149-152°	5g	35.60
met GBC base		
161.21	5g	38.50
	25g	141.00
178.26 mp 185-189°	5g	53.70
cinol page 1123		
H ₂ NC ₆ H ₅ (CH ₃) ₂ (=O)	5g	32.70
483F IRRITANT		
	500mg	96.60
10°C)		
144054-70-6]	1g	103.60
1° (c=1, CHCl ₃)	5g	344.70
FW 136.15	1g	27.30
	10g	151.40
(2-amino-6,7-	500mg	72.80
	1g	121.30



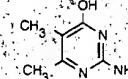
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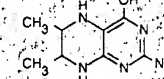
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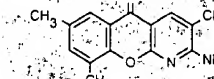
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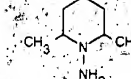
53,461-7



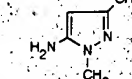
A5,173-4



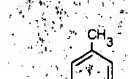
40,244-3



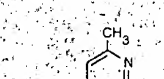
13,345-0



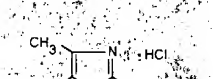
53,222-3



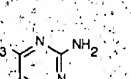
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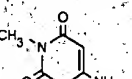
A5,200-5



17,440-8



13,346-9



A5,215-3

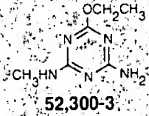
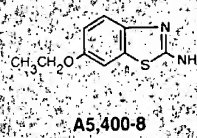
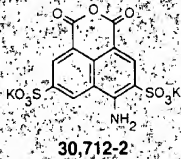
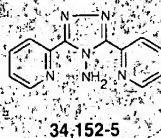
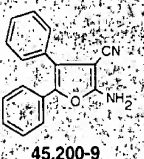
53,461-7	2-Amino-5,6-dimethyl-4-hydroxypyrimidine, 96% [3977-23-9] FW 139.16	1g	89.70
	mp 333-337° IRRITANT		
A5,173-4	2-Amino-6,7-dimethyl-4-hydroxy-5,6,7,8-tetrahydropteridine hydrochloride, 96% [945-43-7] (2-amino-6,7-dimethyl-5,6,7,8-tetrahydro-4-pteridinol hydrochloride) FW 231.69 mp 244-247°(dec.) FT-NMR 1(3),385B FT-IR 1(2),830C Safety 2,164D R&S 1(2),2581E IRRITANT	100mg 250mg	57.80 115.50
	2-Amino-6,7-dimethyl-4-oxo-4H-1-benzopyran-3-carboxaldehyde, see 40,209-5, 2-Amino-3-formyl-6,7-dimethylchromone page 86		
40,244-3	2-Amino-7,9-dimethyl-5-oxo-5H-[1]benzopyrano[2,3-b]pyridine-3-carbonitrile, 98% [206658-82-4] FW 265.27 mp >300° IRRITANT	1g	31.70
12,649-7	4-Amino-2,5-dimethylphenol, tech. [3096-71-7] (4-amino-2,5-xenol) H ₂ NC ₆ H ₂ (CH ₃) ₂ OH FW 137.18 mp 242-243° Beil. 13,634 FT-NMR 1(1),532A FT-IR 1(1),1233B Safety 2,165B R&S 1(1),1425G IRRITANT	5g 10g	85.30 140.40
13,345-0	1-Amino-2,6-dimethylpiperidine, tech., 90% [39135-39-2] FW 128.22 bp 65-80°/30mm n _D 1.4650 d 0.865 Fp 108°F(42°C) FT-NMR 1(1),577B FT-IR 1(1),369B Safety 2,165D R&S 1(1),389H RTECS# TM4175000 IRRITANT	1g 5g	45.30 135.60
	2-Amino-6,7-dimethyl-4-pteridinol, see A5,160-2, 2-Amino-6,7-dimethyl-4-hydroxypteridine page 80		
53,222-3	5-Amino-1,3-dimethylpyrazole, 97% [3524-32-1] FW 111.15 mp 65-69° IRRITANT	10g	125.00
A5,180-7	2-Amino-4,6-dimethylpyridine, 99% [5407-87-4] (6-amino-2,4-lutidine) FW 122.17 mp 63-64° bp 235° FT-NMR 1(3),300C FT-IR 1(2),773A Safety 2,166A R&S 1(2),2529G RTECS# US1818000 IRRITANT	5g 25g	18.90 62.40
A5,200-5	2-Amino-4,6-dimethylpyrimidine, 95% [767-15-7] FW 123.16 mp 151-153° FT-IR 1(2),822D R&S 1(2),2575E IRRITANT	25g 100g	22.60 72.90
	4-Amino-N-(2,6-dimethyl-4-pyrimidinyl)benzenesulfonamide, see 28,721-0, Sulfisomidine page 1700		
	2-Amino-6,7-dimethyl-5,6,7,8-tetrahydro-4-pteridinol hydrochloride, see A5,173-4, 2-Amino-6,7-dimethyl-4-hydroxy-5,6,7,8-tetrahydropteridine hydrochloride, page 81		
17,440-8	2-Amino-4,5-dimethylthiazole hydrochloride, 98% [71574-33-9] FW 164.66 mp 273-274°(dec.) FT-IR 1(2),645B R&S 1(2),2387I	10g 50g	36.70 119.10
13,346-9	3-Amino-5,6-dimethyl-1,2,4-triazine, 97% [17584-12-2] FW 124.15 mp 210-212° FT-NMR 1(3),413C FT-IR 1(2),847A R&S 1(2),2599F RTECS# XY3177100 IRRITANT	25g	50.00
A5,215-3	6-Amino-1,3-dimethyluracil, 98% [6642-31-5] FW 155.16 mp 295°(dec.) Beil. 24,471 FT-NMR 1(1),1321C FT-IR 1(1),812B R&S 1(1),955B RTECS# YQ8755000	100g	40.30
23,874-0	4-Amino-3,5-dinitrobenzamide, 98% [54321-79-8] H ₂ NC ₆ H ₂ (NO ₂) ₂ CONH ₂ FW 226.15 mp 251-254° Beil. 14(2),273 FT-IR 1(2),374C Safety 2,166B R&S 1(2),2023H IRRITANT	5g 25g	10.00 52.60
55,952-0	4-Amino-3,5-dinitrobenzotrifluoride, 97% [445-66-9] H ₂ NC ₆ H ₂ (NO ₂) ₂ CF ₃ FW 251.12 mp 144-149° IRRITANT	5g 25g	40.40 161.60
52,162-0	2-Amino-3,5-dinitrothiophene, 95% [2045-70-7] FW 189.15 mp 174-176° IRRITANT	5g	52.30
	5-Amino-2,6-dioxo-1,2,3,6-tetrahydro-4-pyrimidinecarboxylic acid, see 19,121-3, 5-Aminoorotic acid page 101		
	Aminodiphenylamine, see N-Phenylphenylenediamine		
	4-Aminodiphenylamine diazonium sulfate, see 20,146-4, Variamine Blue RT salt page 1910		
	1-Amino-1,2-diphenylethane, see 13,702-2, 1,2-Diphenylethylamine page 776		
52,370-4	(1R,2R)-(+)-2-Amino-1,2-diphenylethanol, 97% [88082-66-0] C ₆ H ₅ CH(NH ₂)CH(C ₆ H ₅)OH FW 213.28 mp 143-147° [α] _D +7° (c=0.6, C ₂ H ₅ OH)	1g 5g	82.50 274.70

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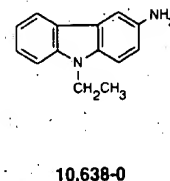
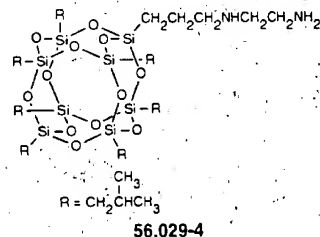
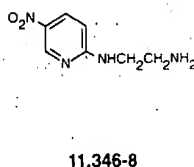
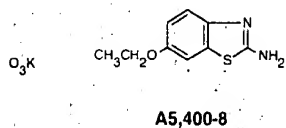
■ Aminodiphe ■

		US \$
33,189-9	(1R,2S)-(-)-2-Amino-1,2-diphenylethanol, 99% [23190-16-1] <chem>C6H5CH(NH2)CH(C6H5)OH</chem> FW 213.28 mp 142-144° [α] _D ²⁵ -7.0° (c=0.6, C ₂ H ₅ OH) Beil 13(3):424 Merck Index 13.449 FT-NMR 1(2):570B R&S 1(1):1465F IRRITANT	1g 28.80 5g 121.30
33,188-0	(1S,2R)-(+)-2-Amino-1,2-diphenylethanol, 99% [23364-44-5] <chem>C6H5CH(NH2)CH(C6H5)OH</chem> FW 213.28 mp 142-144° [α] _D ²⁵ +7.0° (c=0.6, C ₂ H ₅ OH) Beil 13(2):424 Merck Index 13.449 FT-NMR 1(2):570A R&S 1(1):1465E IRRITANT	1g 27.20 5g 113.80
52,371-2	(1S,2S)-(-)-2-Amino-1,2-diphenylethanol, 97% [23190-17-2] <chem>C6H5CH(NH2)CH(C6H5)OH</chem> FW 213.28 mp 140-144° [α] _D ²⁵ -7° (c=0.6, C ₂ H ₅ OH) IRRITANT	1g 73.00 5g 242.90
45,200-9	2-Amino-4,5-diphenyl-3-furancarboxitrile, 95% [5503-73-1] FW 260.30 mp 201-205° Beil 18(5):12,274	1g 25.80 10g 142.00
A5,360-5	Aminodiphenylmethane, 97% [91-00-9] (benzhydramine) (C ₆ H ₅) ₂ CHNH ₂ FW 183.25 mp 12° bp 295° n _D ²⁰ 1.5950 d ₄ ²⁰ 1.063 Fp >230°F (110°C) Beil 12:1323 Merck Index 13.1076 FT-NMR 1(2):568A FT-IR 1(1):1268C Safety 2.166C R&S 1(1):1463N TOXIC IRRITANT	5g 13.70 25g 43.70 100g 123.10
17,688-5	Aminodiphenylmethane hydrochloride, 97% [5267-34-5] (benzhydramine hydrochloride) (C ₆ H ₅) ₂ CHNH ₂ ·HCl FW 219.72 mp 293-295° Beil 12:1323 Merck Index 13.1085 FT-NMR 1(2):568B FT-IR 1(1):1268D Safety 2.166D R&S 1(1):1463O IRRITANT	5g 15.40 25g 53.80 100g 149.40
27,306-6	4-Amino-2,6-diphenylphenol, 98% [50432-01-4] H ₂ NC ₆ H ₂ (C ₆ H ₅) ₂ OH FW 261.32 mp 148-150° Beil 13:735 FT-NMR 1(2):552A R&S 1(1):1445A IRRITANT	250mg 26.90 1g 75.90
54,946-0	(R)-(+)-2-Amino-1,1-diphenyl-1-propanol, 99% [78603-93-7] FW 227.31 mp 102-105° [α] _D ²⁵ +90° (c=1, CHCl ₃) IRRITANT HYGROSCOPIC Product of Onyx Scientific, U.K.	1g 103.60 5g 344.70
38,238-8	(S)-(-)-2-Amino-1,1-diphenyl-1-propanol, 99% [78603-91-5] <chem>CH3CH(NH2)C(C6H5)2OH</chem> FW 227.31 mp 100-102° [α] _D ²⁵ -85° (c=1, CHCl ₃) FT-NMR 1(2):581A R&S 1(1):1471K IRRITANT	100mg 32.50 500mg 124.40
34,152-5	4-Amino-3,5-di-2-pyridyl-4H-1,2,4-triazole, 97% [1671-88-1] FW 238.25 mp 182-184° FT-NMR 1(3):353A R&S 1(2):2559I IRRITANT	1g 34.40 5g 114.50
30,712-2	4-Amino-3,6-disulfo-1,8-naphthalic anhydride, dipotassium salt [79539-35-8] (Lucifer Yellow anhydride, dipotassium salt) FW 449.51 mp >300° Safety 2.167A R&S 1(2):1961E IRRITANT	100mg 170.80
15,924-7	12-Aminododecanoic acid, 95% [693-57-2] H ₂ N(CH ₂) ₁₁ CO ₂ H FW 215.34 mp 185-187° FT-IR 1(1):579C Safety 2.167B R&S 1(1):659K IRRITANT	5g 8.50 25g 33.50
	2-Aminoethanesulfonic acid, see 15,224-2, Taurine, page 1710	
12,292-0	2-Aminoethanethiol hydrochloride, 98% [156-57-0] (cysteamine hydrochloride) H ₂ NCH ₂ CH ₂ SH·HCl FW 113.61 mp 66-68° Beil 4:286 Merck Index 13.2785 FT-IR 1(1):350D Safety 2.167D R&S 1(1):375B RTECS# KJ0200000 HYGROSCOPIC	25g 26.40 100g 87.20
	2-Aminoethanol, see Ethanolamine	
56,299-8	2-Aminoethanol-2-penten-4-one, 97% [312933-35-0] <chem>HOCH2CH2NHC(CH3)=CHCOCH3</chem> FW 143.18 mp 74-78° IRRITANT	1g 39.40
A5,400-8	2-Amino-6-ethoxybenzothiazole, tech. [94-45-1] FW 194.26 mp 161-163° FT-NMR 1(3):208C FT-IR 1(2):704B IRRITANT	25g 18.60 100g 43.80
A5,405-9	2-(2-Aminoethoxy)ethanol, 98% [929-06-6] (diethylene glycolamine) <chem>H2NCH2CH2OCH2CH2OH</chem> FW 105.14 bp 218-224° d ₄ ²⁰ 1.048 Fp >230°F (110°C) Beil 4(3):642 FT-NMR 1(1):542C FT-IR 1(1):339C Safety 2.168A R&S 1(1):365J RTECS# KJ6125000 CORROSIVE	5g 17.90 100g 20.80
52,300-3	2-Amino-4-ethoxy-6-(methylamino)-1,3,5-triazine, 97% [62096-63-3] FW 169.19 mp 171-174° IRRITANT	25g 138.80

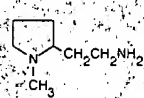


	US \$
1g	28.80
5g	121.30
≈0.6, C ₂ H ₅ OH)	
465F IRRITANT	
1g	27.20
5g	113.80
≈0.6, C ₂ H ₅ OH)	
465E IRRITANT	
1g	73.00
5g	242.90
6, C ₂ H ₅ OH)	
1g	25.80
10g	142.00
C ₆ H ₅) ₂ CHNH ₂	5g 13.70
Beil. 12,1323, Merck	25g 43.70
R&S 1(1),1463N	100g 123.10
inhydramine	5g 15.40
12,1323 Merck	25g 53.80
R&S 1(1),1463O	100g 149.40
15)OH FW 261.32	250mg 26.90
IRITANT	1g 75.90
W 227.31	1g 103.60
	5g 344.70
100mg	32.50
500mg	124.40
CHCl ₃)	
W 238.25	1g 34.40
	5g 114.50
salt [79539-35-8]	100mg 170.80
300° Safety 2,167A	
FW 215.34	5g 8.50
IRITANT	25g 33.50
'10	
nine	25g 26.40
4,286 Merck	100g 87.20
EC5# KJ0200000	
1g	39.40
VT	
ip 161-163°	25g 18.60
	100g 43.80
amine)	5g 17.90
30°F(110°C)	100g 20.80
R&S 1(1),365J	
16-63-3] FW 169.19	25g 138.80

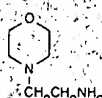
35,962-9	L-α-[2-(2-Aminoethoxy)vinyl]glycine hydrochloride, 95% [55720-26-8] 5mg 106.70
	H ₂ NCH ₂ CH ₂ OCH=CHCH(NH ₂)CO ₂ H · HCl FW 196.64 [α] _D ²⁵ +83° (c=1, phosphate buffer) 25mg 372.30
	R&S 1(1),669H RTECS# EM9080000 TOXIC
	Contains 2-3% D-isomer
	N-(2-Aminoethyl)acetamide, see 39,726-1, N-Acetylenediamine page 23
12,758-2	2-(2-Aminoethylamino)ethanol, 99% [111-41-1] H ₂ NCH ₂ CH ₂ NHCH ₂ CH ₂ OH 100g 21.00
★	FW 104.15 bp 238-240°/752mm n _D ²⁰ 1.4860 d 1.030 Fp >230°F(110°C) Beil. 4,286 1kg 32.70
	FT-NMR 1(1),552C FT-IR 1(1),347D Safety 2,168B R&S 1(1),371H 3kg 83.60
	RTECS# KJ6300000 CORROSIVE LACHRYMATOR
11,346-8	2-(2-Aminoethylamino)-5-nitropyridine, 96% [29602-39-9] FW 182.18 10g 53.30
	mp 126-128° FT-IR 1(2),776D Safety 2,169C R&S 1(2),2531O IRRITANT 25g 87.50
	Aminoethylaminopropylisobutyl-POSS, see 56,029-4, 1-[3-(2-Aminoethyl)-
	amino]propyl-3,5,7,9,11,13,15-isobutylpentacyclo-[9.5.1.1(3.9).1(5,15).1(7,13)]-
	octasiloxane page 83
56,029-4	1-[3-(2-Aminoethylamino)propyl-3,5,7,9,11,13,15-isobutyl- 1g 26.70
NEW	pentacyclo[9.5.1.1(3.9).1(5,15).1(7,13)]octasiloxane(aminoethylaminopropyl- 5g 88.90
	isobutyl-POSS) FW 917.65 mp 109-116°
	[3-(2-Aminoethylamino)propyl]trimethoxysilane, see N-[3-(Trimethoxysilyl)-
	propyl]ethylenediamine
27,524-7	4-(2-Aminoethyl)benzenesulfonamide, 99% [35303-76-5] 25g 28.20
	H ₂ NCH ₂ CH ₂ C ₆ H ₄ SO ₂ NH ₂ FW 200.26 mp 150-152° Beil. 14(3),2234 100g 78.20
	FT-NMR 1(2),1637C R&S 1(2),2243E
53,841-8	4-(2-Aminoethyl)benzoic acid, 95% [1199-69-5] H ₂ NCH ₂ CH ₂ C ₆ H ₄ CO ₂ H 5g 24.40
NEW	FW 165.19 mp >350° 25g 81.00
	α-(1-Aminoethyl)benzyl alcohol, see Norephedrine
10,638-0	3-Amino-9-ethylcarbazole, tech., 90% [132-32-1] FW 210.28 mp 98-100° 25g 19.90
★	Beil. 22(1),642 FT-NMR 1(3),163A FT-IR 1(2),680D Safety 2,169D R&S 1(2),2425I
	RTECS# FE3590000 CANCER SUSPECT AGENT: MUTAGEN
	Indicator for peroxidase activity. Anal. Biochem. 1973, 56, 353. Am. J. Clin. Pathol.
	1975, 63, 451.
	May contain up to 3% ethanol
	2-Aminoethyl dibutylborinate, see 44,640-8, Dibutylborinic acid, ethanolamine
	ester page 594
29,286-9	2-Aminoethyl dihydrogenphosphate, 98% [1071-23-4] H ₂ NCH ₂ CH ₂ OP(O)(OH) ₂ 5g 14.00
★	FW 141.07 mp 241-243° Beil. 4(3),644 Safety 2,170A R&S 1(1),1115B 25g 46.40
	RTECS# KJ5830000 CORROSIVE
	α-(1-Aminoethyl)-3,4-dihydroxybenzyl alcohol, see 37,551-9, (-)-3,4-Dihydroxy-
	norephedrine page 685
	2-Aminoethyl diphenylborinate, see 12,670-5, Diphenylborinic acid, ethanol-
	amine ester page 775
	2-Aminoethyl ether dihydrochloride, see 17,609-5, 2,2'-Oxybis(ethylamine)
	dihydrochloride page 1398
A5,440-7	2-Aminoethyl hydrogensulfate, 97% [926-39-6] H ₂ NCH ₂ CH ₂ OSO ₃ H FW 141.15 25g 12.60
★	mp 277°(dec.) Beil. 4,276 FT-IR 1(1),904B Safety 2,170B R&S 1(1),1079J 100g 34.90
	RTECS# KJ6371000 IRRITANT 500g 82.60
A5,445-8	α-(1-Aminoethyl)-4-hydroxybenzyl alcohol hydrochloride, 99+% [35999-22-5] 10g 44.80
	(4-hydroxynorephedrinehydrochloride) HOC ₆ H ₄ CH(CH(NH ₂))CH ₂ OH · HCl
	FW 203.67 mp 202°(dec.) [α] _D ²² 0° (c=1, H ₂ O) FT-NMR 1(2),613A FT-IR 1(1),1290B
	Safety 2,170C R&S 1(1),1489L RTECS# DN4900000 IRRITANT
	3-(2-Aminoethyl)indole, see Tryptamine



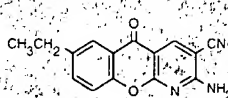
A5,460-1	S-(2-Aminoethyl)isothiuronium bromide hydrobromide, 99% [56-10-0].....	25g	US \$ 35.40
★	[2-(2-aminoethyl)-2-thiopseudouréa dihydrobromide] H ₂ NC(SCH ₂ CH ₂ NH ₂)=NH·2HBr FW 281.02 mp 190-196° Merck Index 13,178 FT-IR 1(1),816C Safety 2,170D R&S 1(1),959N RTECS# UM0175000 MOISTURE-SENSITIVE TOXIC Has radioprotective activity against X-rays. Indian J. Exp. Biol. 1993, 31, 837, ibid. 1992, 30, 478.	100g	104.20
51,615-5	2-Aminoethyl methacrylate hydrochloride, 90% [2420-94-2].....	5g	45.90
★	H ₂ C=C(CH ₃)CO ₂ CH ₂ CH ₂ NH ₂ ·HCl FW 165.62 mp 102-110° IRRITANT	25g	152.40
	3-(2-Aminoethyl)-5-methylindole hydrochloride, see 13,422-8, 5-Methyl-tryptamine hydrochloride page 1303		
13,950-5	2-(2-Aminoethyl)-1-methylpyrrolidine, 97% [51387-90-7] FW 128.22 n _D 1.4680 d 0.885 Fp 149°F (65°C) FT-NMR 1(1),562B FT-IR 1(1),368D Safety 2,171B R&S 1(1),379O MOISTURE-SENSITIVE IRRITANT	1g	26.90
		5g	88.10
A5,500-4	4-(2-Aminoethyl)morpholine, 99% [2038-03-1] FW 130.19 bp 205° n _D 1.4750 d 0.992 Fp 347°F (175°C) Beil. 27(4),370 FT-NMR 1(1),600C FT-IR 1(1),384C Safety 2,171C R&S 1(1),403J RTECS# QD7350000 CORROSIVE TOXIC	5g	12.10
★		25g	40.10
		100g	96.80
40,228-1	2-Amino-7-ethyl-5-oxo-5H-[1]benzopyrano[2,3-b]pyridine-3-carbonitrile, 98% [68302-09-0] FW 265.27 mp >300° IRRITANT	1g	29.80
	4-(2-Aminoethyl)phenol, see Tyramine		
49,672-3	4-(1-Aminoethyl)phenol propoxylate CH ₃ CH(NH ₂)C ₆ H ₄ (OC ₃ H ₇) ₂ OH n _D 1.4600 d 1.016 Fp >230°F (110°C) IRRITANT Average M _n ca. 1.880	20mL	19.10
36,754-0	(R)-(-)-(1-Aminoethyl)phosphonic acid, 99% [60687-36-7] CH ₃ CH(NH ₂)P(O)(OH) ₂ FW 125.07 mp 290° (dec.) [α] _D -4.8° (c=5, H ₂ O) FT-NMR 1(1),1476A R&S 1(1),1101J IRRITANT	100mg	29.20
		1g	160.90
36,755-9	(S)-(+)-(1-Aminoethyl)phosphonic acid, 99% [66068-76-6] CH ₃ CH(NH ₂)P(O)(OH) ₂ FW 125.07 mp 290° (dec.) [α] _D +4.8° (c=5, H ₂ O) FT-NMR 1(1),1476B R&S 1(1),1101K IRRITANT	100mg	26.80
		1g	146.30
26,867-4	2-Aminoethylphosphonic acid, 99% [2041-14-7] H ₂ NCH ₂ CH ₂ P(O)(OH) ₂ FW 125.07 mp 296° (dec.) Beil. 4(3),1792 FT-NMR 1(1),1477A Safety 2,172A R&S 1(1),1101M IRRITANT	250mg	29.30
		1g	84.90
A5,520-9	1-(2-Aminoethyl)piperazine, 99% [140-31-8] FW 129.21 bp 218-222° n _D 1.5000 d 0.985 Fp 200°F (93°C) FT-NMR 1(1),586A FT-IR 1(1),374D Safety 2,172B R&S 1(1),395A RTECS# TK8050000 CORROSIVE TOXIC Utilized in a variety of reactions for studying corrosion inhibition, biological activity ² and metal ligand effects on catalysis. ³ (1) J. Chem. Soc., Perkin Trans. 2 1992, 939; (2) J. Med. Chem. 1993, 36, 765; (3) Tetrahedron 1992, 48, 1999.	5g	15.40
★		100g	19.20
		500g	39.20
14,166-6	1-(2-Aminoethyl)piperidine, 98% [27578-60-5] FW 128.22 bp 186° n _D 1.4740 d 0.899 Fp 136°F (57°C) Beil. 20,67 FT-NMR 1(1),575C FT-IR 1(1),369D Safety 2,172C R&S 1(1),389B CORROSIVE	5g	29.60
		10g	47.80
12,715-9	N-(2-Aminoethyl)-1,3-propanediamine, 97% [13531-52-7] H ₂ N(CH ₂) ₃ NHCH ₂ CH ₂ NH ₂ FW 117.20 n _D 1.4810 d 0.928 Fp 205°F (96°C) FT-NMR 1(1),500A FT-IR 1(1),308D Safety 2,173A R&S 1(1),331C CORROSIVE	25g	22.40
★		100g	62.40
29,576-0	5-Amino-1-ethylpyrazole, 99% [3528-58-3] FW 111.15 mp 52-57° bp 92-94°/0.01mm Fp >230°F (110°C) FT-NMR 1(3),70B Safety 2,173C R&S 1(2),2347B CORROSIVE HYGROSCOPIC Heterocyclic building block. J. Chem. Soc., Perkin Trans. 1 1980, 938.	1g	40.60
		5g	161.60
A5,530-6	2-(2-Aminoethyl)pyridine, 95% [2706-56-7] FW 122.17 bp 92-93°/12mm n _D 1.5360 d 1.021 Fp 213°F (100°C) Beil. 22,434 FT-NMR 1(3),289C FT-IR 1(2),764C Safety 2,173D R&S 1(2),2521H IRRITANT	1g	15.40
★		10g	82.70
		50g	310.30
51,811-5	3-(2-Aminoethyl)pyridine dihydrobromide, 98% [307496-23-7] FW 284.01 mp 248-252° IRRITANT	1g	33.20
		5g	109.40



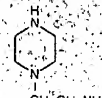
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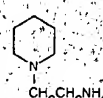
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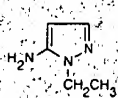
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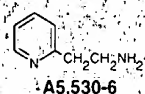
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14,166-6



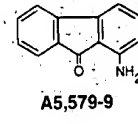
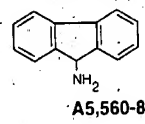
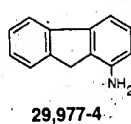
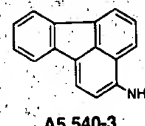
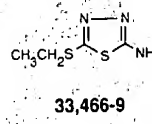
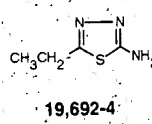
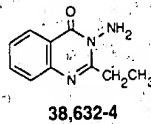
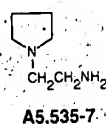
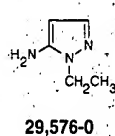
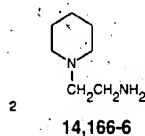
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A5,530-6

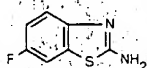
		US \$
10-0	25g	35.40
3,178	100g	104.20
ibid. 1992,		
	5g	45.90
	25g	152.40
ethyl-		
n _D 1.4680	1g	26.90
18	5g	88.10
n _D 1.4750	5g	12.10
184C	25g	40.10
C	100g	96.80
nitrile, 98%	1g	29.80
n _D 1.4600	20mL	19.10
(H ₂ P(O)(OH) ₂	100mg	29.20
S 1(1), 1101J	1g	160.90
	100mg	26.80
	1g	146.30
OH ₂	250mg	29.30
172A	1g	84.90
n _D 1.5000	5g	15.40
72B	100g	19.20
	500g	39.20
activity ² and		
939. (2) J.		
n _D 1.4740	5g	29.60
Safety 2, 172C	10g	47.80
	25g	22.40
C)	100g	62.40
ROSIVE		
	1g	40.60
	5g	161.60
mm n _D 1.5360	1g	15.40
4C	10g	82.70
	50g	310.30
V 284.01	1g	33.20
	5g	109.40

A5,535-7	1-(2-Aminoethyl)pyrrolidine, 98% [7154-73-6] FW 114.19 bp 66-70°/23mm	1g	11.60
★	n _D 1.4680 d 0.901 Fp 118°F(47°C) FT-NMR 1(1), 561C FT-IR 1(1), 358A Safety 2, 174B	5g	38.10
	R&S 1(1), 379M CORROSIVE MOISTURE-SENSITIVE		
38,632-4	3-Amino-2-ethyl-4(3H)-quinazolinone, 99% [50547-51-8] FW 189.22 mp 121-123°	1g	36.80
	R&S 1(2), 2647E IRRITANT	5g	124.00
	The N-acetoxy derivative is used for azidination of alkenes ¹ and reacts with vinyl-silanes and -stannanes to yield silyl or stannyl-substituted azidines. ² (1) Tetrahedron 1989, 45, 2875. (2) J. Chem. Soc., Chem. Commun. 1989, 836.		
32,919-3	2-Amino-4-(ethylsulfonyl)phenol, 97% [43115-40-8] H ₂ NC ₆ H ₃ (SO ₂ C ₂ H ₅)OH	250g	18.70
	FW 201.25 mp 128-131° FT-NMR 1(2), 1586C R&S 1(2), 2187N IRRITANT	1kg	50.80
19,692-4	2-Amino-5-ethyl-1,3,4-thiadiazole, 97% [14068-53-2] FW 129.18 mp 200-203°	5g	16.60
★	FT-NMR 1(3), 118A FT-IR 1(2), 651A R&S 1(2), 2393G RTECS# X13095000 IRRITANT	25g	53.40
	2-Amino-4-(ethylthio)butyric acid, see Ethionine		
27,780-0	S-(2-Aminoethyl) thiophosphate, monosodium salt, 97% [3724-89-8]	10mg	19.50
	H ₂ NCH ₂ CH ₂ SP(O)(OH)ONa FW 179.11 mp 220°(dec.) Beil. 4(4), 1578 Merck	50mg	73.00
	Index 13,7427 R&S 1(1), 1113A RTECS# TE6483500 IRRITANT		
	2-(2-Aminoethyl)-2-thiopseudourea dihydrobromide, see A5,460-1, S-(2-Aminoethyl)isothiuronium bromide hydrobromide page 84		
33,466-9	2-Amino-5-(ethylthio)-1,3,4-thiadiazole, 99% [25660-70-2] FW 161.25	5g	18.20
★	mp 135-137° Beil. 27(2), 785 FT-NMR 1(3), 118C R&S 1(2), 2393J	25g	67.70
28,455-6	(2-Aminoethyl)trimethylammonium chloride hydrochloride, 99% [3399-67-5]	1g	50.50
	H ₂ NCH ₂ CH ₂ N(CH ₃) ₃ Cl·HCl FW 175.10 mp 260°(dec.) Beil. 4(2), 690 Safety 2, 174C	5g	167.90
	R&S 1(1), 423E IRRITANT HYGROSCOPIC		
A5,540-3	3-Amino-9-fluorenone, 90% [2693-46-1] (3-fluoranthraniline) FW 217.27	250mg	22.60
	mp 115-117° FT-IR 1(1), 1256A Safety 2, 174D R&S 1(1), 1451D RTECS# LL4000000	1g	62.50
		10g	346.70
		100mg	43.00
29,977-4	1-Amino-9-fluorenone, 99% [6344-63-4] (1-fluorenamine) FW 181.24 mp 125-126°		
	Beil. 12(3), 3285 Safety 2, 175A R&S 1(1), 1449B RTECS# LL5074000 IRRITANT	5g	35.00
A5,550-0	2-Amino-9-fluorenone, 98% [153-78-6] (2-fluorenamine) FW 181.24 mp 127-131°	25g	133.70
	Beil. 12, 1331 FT-IR 1(1), 1254C Safety 2, 175B R&S 1(1), 1449C RTECS# LL5075000	5g	40.00
A5,560-8	8-Amino-9-fluorenone hydrochloride, 98% [5978-75-6] (9-fluorenamine hydrochloride) FW 217.70 mp 265-270°(dec.) Beil. 12, 1331 FT-NMR 1(2), 634B	25g	158.10
	FT-IR 1(1), 1301A R&S 1(1), 1503L		
	2-Amino-9-fluorenone, see A5,680-9, 2-Amino-9-hydroxyfluorenone page 88		
A5,579-9	1-Amino-9-fluorenone, 97% [6344-62-3] FW 195.22 mp 118-120° Beil. 14, 113	250mg	35.80
	FT-NMR 1(2), 906A FT-IR 1(2), 76B Safety 2, 175D R&S 1(2), 1679J IRRITANT	1g	120.00
A5,580-2	2-Amino-9-fluorenone [3096-57-9] FW 195.22 mp 157-160° Beil. 14, 113	1g	37.70
	FT-NMR 1(2), 906B FT-IR 1(2), 76C Safety 2, 176A R&S 1(2), 1679K RTECS# LL8980200	10g	208.80
	IRRITANT		
18,922-7	3-Amino-9-fluorenone, 98% [6276-05-7] FW 195.22 mp 154-155° Beil. 14(2), 69	100mg	51.90
	Safety 2, 176B R&S 1(2), 1679N IRRITANT		
12,294-7	4-Amino-9-fluorenone, 95% [4269-15-2] FW 195.22 mp 138-140° Beil. 14, 113	1g	140.40
	FT-NMR 1(2), 906C FT-IR 1(2), 77A Safety 2, 176C R&S 1(2), 1679O		
	RTECS# LL8980300 IRRITANT		
	Amino-9-fluorescein, see Fluoresceinamine		
37,016-9	2-Amino-4-fluorobenzoic acid, 97% [446-32-2] (4-fluoroanthranilic acid)	250mg	12.00
	H ₂ NC ₆ H ₃ (F)CO ₂ H FW 155.13 mp 194-196° Beil. 14(3), 952 FT-NMR 1(2), 1118B	1g	33.00
	R&S 1(2), 1807F IRRITANT		
36,798-2	2-Amino-5-fluorobenzoic acid, 97% [446-08-2] (5-fluoroanthranilic acid)	250mg	15.50
	H ₂ NC ₆ H ₃ (F)CO ₂ H FW 155.13 mp 182-184° Beil. 14(4), 1071 FT-NMR 1(2), 1118C	1g	39.90
	R&S 1(2), 1807G IRRITANT		
	Used in the synthesis of styrylquinazolinones which are potential anticancer agents. J. Med. Chem. 1990, 33, 1721.		

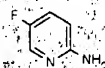


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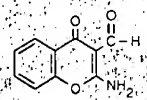
			US \$
44,302-6	2-Amino-6-fluorobenzoic acid, 98% [434-76-4] (6-fluoroanthranilic acid) <chem>Nc1ccc(F)c(C(=O)O)c1</chem> FW 155.13 mp 167-169° IRRITANT. Recently used in the synthesis of a novel benzamide with potent neuroleptic activity. <i>J. Med. Chem.</i> 1996, 39, 1172.	1g 5g	27.20 99.00
42,985-6	2-Amino-6-fluorobenzonitrile, 99% [77326-36-4] (6-fluoroanthranilonitrile) <chem>N#Cc1ccc(F)c(N)c1</chem> FW 136.13 mp 125-128° IRRITANT A medicinal chemistry intermediate used for the preparation of tacrine-related compounds ¹ and antilobate and antibacterial quinazoline derivatives. ² (1) <i>Eur. J. Med.</i> <i>Chem.</i> 1994, 29, 205. (2) <i>J. Med. Chem.</i> 1990, 33, 434.	1g 10g	13.80 76.10
32,421-3	2-Amino-6-fluorobenzothiazole, 99% [348-40-3] FW 168.19 mp 183-185° <chem>Nc1ccc(F)c2c(N1)sc3ccccc32</chem> FW 168.19 mp 183-185° <i>Beil.</i> 27(4), 4858 FT-NMR 1(3), 207A R&S 1(2), 2457A IRRITANT Aminofluorobenzotrifluoride; see Fluoro(trifluoromethyl)aniline	1g 5g	18.50 61.10
54,183-4	2-Amino-6-fluorobenzylamine, 97% [175277-93-7] <chem>Nc1ccc(F)c(CCN)c1</chem> FW 140.16 bp 78-80°/0.8mm n _D 1.5700 d 1.191 Fp >230°F (110°C)	1g	52.80
55,109-0	(R)-(+)-2-Amino-1-fluoro-3-methyl-1,1-diphenylbutane, 97% <chem>CC(C1=CC=CC=C1)(C(F)C2=CC=CC=C2)C(N)=CC3=CC=CC=C3</chem> FW 257.35 mp 54-57° [α] _D +45° (c=1, CHCl ₃) IRRITANT A Product of Onyx Scientific, U.K.	500mg	121.50
55,107-4	(S)-(-)-2-Amino-1-fluoro-3-methyl-1,1-diphenylbutane, 97% <chem>CC(C1=CC=CC=C1)(C(F)C2=CC=CC=C2)C(N)=CC3=CC=CC=C3</chem> FW 257.35 mp 54-57° [α] _D -45° (c=1, CHCl ₃) IRRITANT A Product of Onyx Scientific, U.K.	500mg	121.50
55,463-4	(R)-(+)-2-Amino-1-fluoro-4-methyl-1,1-diphenylpentane, 98% <chem>CC(C1=CC=CC=C1)(C(F)C2=CC=CC=C2)C(N)=CC3=CC=CC=C3</chem> FW 271.37 mp 85-89° [α] _D +51° (c=1, CHCl ₃) IRRITANT 99% ee/HPLC	500mg	121.70
55,464-2	(S)-(-)-2-Amino-1-fluoro-4-methyl-1,1-diphenylpentane, 98% [274674-22-5] <chem>CC(C1=CC=CC=C1)(C(F)C2=CC=CC=C2)C(N)=CC3=CC=CC=C3</chem> FW 271.37 mp 86-89° [α] _D -50° (c=1, CHCl ₃) IRRITANT 99% ee/HPLC	500mg	121.70
51,868-9	2-Amino-5-fluoropyridine, 97% [21717-96-4] FW 112.11 mp 93-97° <chem>Nc1cc(F)ccn1</chem> FW 112.11 mp 93-97° <i>Beil.</i> 22(5), 8,449 IRRITANT 4-Aminofolic acid, see 86,227-4; Aminopterin page 107	250mg	69.20
40,202-8	2-Amino-3-formylchromone, 97% [61424-76-8] (2-amino-4-oxo-4H-1-benzo- pyran-3-carboxaldehyde) FW 189.17 mp 249° (dec.) <i>Beil.</i> 18(5), 11,571 IRRITANT	1g	24.20
40,209-5	2-Amino-3-formyl-6,7-dimethylchromone [94978-87-7] (2-amino-6,7-dimethyl- 4-oxo-4H-1-benzopyran-3-carboxaldehyde) FW 217.22 mp >300° IRRITANT Amino-G acid, see 14,644-7, 7-Amino-2,7-naphthalenedisulfonic acid, monopotassium salt page 98	1g	32.00
25,919-5	Aminoglutethimide, 99% [125-84-8] 3-(4-aminophenyl)-3-ethyl-2,6-piperidine- dione] FW 232.29 mp 152-154° [α] _D 25.0° (c=1, CH ₃ OH) <i>Beil.</i> 22(4), 6596 Merck <i>Index</i> 13,439 FT-NMR 1(2), 1456C FT-IR 1(2), 404A Safety 2,177A R&S 1(2), 2079F RTECS# MA4026950 TERATOGEN IRRITANT	100mg 500mg	34.90 112.00
54,587-2	(R)-(+)-Aminoglutethimide, 97% [55511-44-9] FW 232.29 mp 115.5-119.5° [α] _D +160° (c=1, CH ₃ OH) TERATOGEN IRRITANT	500mg	79.30
54,586-4	(S)-(-)-Aminoglutethimide, 97% [57288-03-6] FW 232.29 mp 115.5-119.5° [α] _D -160° (c=1, CH ₃ OH) TERATOGEN IRRITANT	500mg	79.30
10,926-6	Aminoguanidine bicarbonate, 97% [2582-30-7] H ₂ NNHC(=NH)NH ₂ ·H ₂ CO ₃ FW 136.11, mp 170-172° (dec.) <i>Beil.</i> 3, 117 FT-IR 1(1), 820A R&S 1(1), 963M RTECS# FG1772000	5g 100g 500g	10.40 12.90 30.00
39,649-4	Aminoguanidine hydrochloride, 98+% [1937-19-5] H ₂ NNHC(=NH)NH ₂ ·HCl FW 110.55 mp 165-168° <i>Beil.</i> 3, 118 Merck <i>Index</i> 13,440 RTECS# ME8430000	25g 100g	26.00 71.00
A5,610-8	Aminoguanidine nitrate, 99% [10308-82-4] H ₂ NNHC(=NH)NH ₂ ·HNO ₃ FW 137.11 mp 145-147° <i>Beil.</i> 3, 117 FT-IR 1(1), 820B Safety 2,177B R&S 1(1), 963N OXIDIZER IRRITANT	5g 100g	16.10 40.70
A5,620-5	2-Aminoheptane, 99% [123-82-0] (1-methylhexylamine) CH ₃ (CH ₂) ₄ CH(NH ₂)CH ₃ FW 115.22 bp 142-144° n _D 1.4170 d 0.766 Fp 130°F (54°C) <i>Beil.</i> 4, 194 Merck <i>Index</i> 13,9873 FT-NMR 1(1), 461B FT-IR 1(1), 289A Safety 2,177C R&S 1(1), 307G RTECS# MQ5425000 IRRITANT	25g 100g	28.50 87.50



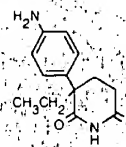
32,421-3



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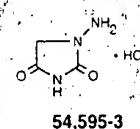
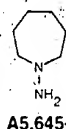
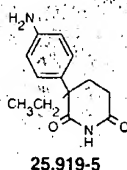


40,202-8



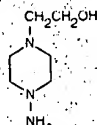
25,919-5

			US \$				US \$
milic acid)	1g	27.20	46,262-4 (R)-(-)-2-Aminoheptane [6240-90-0] CH ₃ (CH ₂) ₄ CH(NH ₂)CH ₃ FW 115.22 bp 142-144°	1g	83.30		US \$
leptic activity.	5g	99.00	nb 1.4190 d 0.766 Fp 130°F(54°C) IRRITANT				
amilonitrile)	1g	13.80	46,263-2 (S)-(+)-2-Aminoheptane [44745-29-1] CH ₃ (CH ₂) ₄ CH(NH ₂)CH ₃ FW 115.22	1g	83.30		
-related	10g	76.10	bp 142-144° nb 1.4190 d 0.766 Fp 130°F(54°C) [α] _D +7° (neat) FLAMMABLE LIQUID				
(1) Eur. J. Med			IRRITANT				
183-185°	1g	18.50	28,463-7 7-Aminoheptanoic acid, 98% [929-17-9] H ₂ N(CH ₂) ₅ CO ₂ H FW 145.20 mp 192-195°	100mg	23.30		
	5g	61.10	Boil 4.459 FT-NMR 1(1).879A Safety 2.178A R&S 1(1).659H RTECS# MJ1770000	1g	153.90		
CH ₃ NH ₂	1g	52.80	IRRITANT	5g	513.20		
CHCl ₃) IRRITANT	500mg	121.50	N-Aminohexamethyleneimine, see A5.645-0. 1-Aminohomopiperidine page 87				
CHCl ₃) IRRITANT	500mg	121.50	2-Aminohexanoic acid, see Norleucine.				
(c=1, CHCl ₃)	500mg	121.70	6-Aminohexanoic acid, see A4.460-6. 6-Aminocaproic acid page 74				
74674-22-5]	500mg	121.70	23,767-1 DL-2-Amino-1-hexanol, tech., 90% [5665-74-7] (DL-norleucinol)	1g	31.70		
(c=1, CHCl ₃)			CH ₃ (CH ₂) ₄ CH(NH ₂)CH ₂ OH FW 117.19 bp 190-192°/740mm nb 1.4520 d 0.912	10g	199.20		
7°	250mg	69.20	Fp 212°F(100°C) Boil. 4(3).809 FT-NMR 1(1).541A FT-IR 1(1).338C Safety 2.178B				
o-4H-1-benzo-	1g	24.20	R&S 1(1).363N IRRITANT HYGROSCOPIC				
571 IRRITANT			53,455-2 (R)-2-Amino-1-hexanol, 97% [80696-29-3] CH ₃ (CH ₂) ₄ CH(NH ₂)CH ₂ OH FW 117.19	1g	62.50		
o-6,7-dimethyl-	1g	32.00	mp 40-44° bp 216-218° Fp 211°F(99°C) [α] _D -14° (c=1, CHCl ₃) IRRITANT				
0° IRRITANT			53,456-0 (S)-2-Amino-1-hexanol, 97% [80696-29-3] (L-norleucinol)	1g	62.50		
acid.			CH ₃ (CH ₂) ₄ CH(NH ₂)CH ₂ OH FW 117.19 mp 35-40° bp 216-218° Fp 211°F(99°C)				
1-2,6-piperidine	100mg	34.90	[α] _D +14° (c=1, CHCl ₃) IRRITANT				
596 Merck	500mg	112.00	A5,635-3 6-Amino-1-hexanol, 97% [4048-33-3] H ₂ N(CH ₂) ₅ OH FW 117.19 mp 56-58°	1g	10.50		
\$S 1(2).2079F			bp 135-140°/30mm Boil. 4(2).748 FT-NMR 1(1).540C FT-IR 1(1).339B Safety 2.178C	5g	35.40		
5-119.5°	500mg	79.30	R&S 1(1).363M RTECS# MO8840000 IRRITANT HYGROSCOPIC	25g	116.20		
5-119.5°	500mg	79.30	23,125-8 1-Aminohexylamide gel [poly(acrylic acid 6-aminohexylamide)]	100g	300.50		
NH ₂ ·H ₂ CO ₃	5g	10.40	[CH ₂ CH[CONH(CH ₂) ₅ NH ₂]] _n Safety 2.178D R&S 1(3).3573A IRRITANT	10g	48.00		
.963M	100g	12.90	HYGROSCOPIC				
500g		30.00	ca. 2.4 meq/g				
NH(NH ₂)·HCl	25g	26.00	47,185-2 (1S)-(+)-(1-Aminohexyl)phosphonic acid, 98% CH ₃ (CH ₂) ₄ CH(NH ₂)P(O)(OH) ₂	250mg	101.70		
E8430000	100g	71.00	FW 181.17 mp 270-276° [α] _D +25° (c=5, 1N NaOH) IRRITANT				
HNO ₃ FW 137.11	5g	16.10	08088 4-Aminohippuric acid, puriss., ≥99.0% (T) [61-78-9] [N-(4-	10g	37.60		
3N OXIDIZER	100g	40.70	aminobenzoyl)glycine]	50g	141.10		
H ₂) ₄ CH(NH ₂)CH ₃	25g	28.50	12,295-5 4-Aminohippuric acid, 97% [61-78-9] H ₂ NC ₆ H ₄ CONHCH ₂ CO ₂ H FW 194.19	25g	34.00		
94 Merck	100g	87.50	mp 199-200°(dec.) Beil. 14(2).258 Merck Index 13.441 FT-NMR 1(2).1408A	100g	114.00		
\$S 1(1).307G			FT-IR 1(2).378B R&S 1(2).2027G IRRITANT				
			A5,645-0 1-Aminohomopiperidine, 95% [5906-35-4] (N-aminohexamethyleneimine)	5g	106.40		
			FW 114.19 bp 165° nb 1.4850 d 0.984 Fp 133°F(56°C) FT-NMR 1(1).579C				
			FT-IR 1(1).376B Safety 2.179B R&S 1(1).391C RTECS# CM3165000 IRRITANT				
			54,595-3 1-Aminohydantoin hydrochloride, 98% [2827-56-7] FW 151.55 mp 201-205°	5g	30.30		
			IRRITANT	25g	100.70		
			4-Amino-3-hydrazino-5-mercapto-1,2,4-triazole, see 16,289-2. Purpald [®]				
			page 1602				
			4-Amino-5-hydrazino-4H-1,2,4-triazole-3-thiol, see 16,289-2. Purpald [®]				
			page 1602				
			β-Aminohydrocinnamic acid, see 15,949-2. 3-Amino-3-phenylpropionic acid				
			page 104				
			38,777-0 3-Amino-4-hydroxybenzenesulfonic acid hydrate [98-37-3]	50g	15.40		
			H ₂ NC ₆ H ₃ (OH)SO ₃ H·xH ₂ O FW 189.19 mp 220°(dec.) Beil. 14.814 R&S 1(2).2197H	250g	51.60		
			2-Amino-3-hydroxybenzoic acid; see 14,877-6. 3-Hydroxyanthranilic acid				
			page 1020				
			28,964-7 3-Amino-4-hydroxybenzoic acid, 97% [1571-72-8] H ₂ NC ₆ H ₃ (OH)CO ₂ H	5g	22.10		
			FW 153.14 mp 208°(dec.) Beil. 14.593 FT-NMR 1(2).1122C Safety 2.180A	25g	95.90		
			R&S 1(2).1809E IRRITANT LIGHT-SENSITIVE				
			33,959-8 4-Amino-3-hydroxybenzoic acid [2374-03-0] H ₂ NC ₆ H ₃ (OH)CO ₂ H FW 153.14	5g	32.60		
			mp 211-215° Beil. 14.589 FT-NMR 1(2).1123A R&S 1(2).1809F IRRITANT	25g	86.70		
			2-Amino-4-hydroxybutyric acid, see Homoserine				

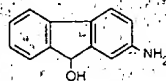


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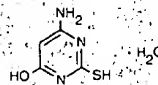
			US\$
46,735-9	(S)-(-)-4-Amino-2-hydroxybutyric acid, 96% [40371-51-5] $H_2NCH_2CH_2CH(OH)CO_2H$ FW 119.12 mp 200-203° [α] _D ²³ -30° (c=1, H ₂ O) Beil. 4(4):3170. IRRITANT	5g 25g	40.80 135.20
A5,665-5	(±)-4-Amino-3-hydroxybutyric acid, 98% [924-49-2] $H_2NCH_2CH(OH)CH_2CO_2H$ FW 119.12 mp 223°(dec.) Beil. 4(2):938 Merck Index 13,456 FT-NMR 1(1):882C FT-IR 1(1):581D R&S 1(1):661K RTECS# ES7015000 IRRITANT	1g 5g	18.80 62.10
	2-Amino-4-hydroxy-6,7-dimethyl-5,6,7,8-tetrahydropteridinehydrochloride, see A5,173-4; 2-Amino-6,7-dimethyl-4-hydroxy-5,6,7,8-tetrahydropteridine hydrochloride page 81		
12,296-3	1-Amino-4-(2-hydroxyethyl)piperazine, 96% [3973-70-4] (4-amino-1-piperazine- ethanol) FW 145.21 mp 105-107° FT-NMR 1(1):585C FT-IR 1(1):374B R&S 1(1):393M IRRITANT	1g 5g	118.80 397.30
A5,680-9	2-Amino-9-hydroxyfluorene, 97% [33417-27-5] (2-amino-9-fluorene) FW 197.24 mp 197-199° Beil. 13,722 Safety 2,180B R&S 1(1):1449L RTECS# LL8410000 IRRITANT	100mg	31.20
A5,740-6	4-Amino-6-hydroxy-2-mercaptopyrimidine monohydrate, 98% [65802-56-4] (6-amino-2-mercapto-4-pyrimidinol-6-amino-2-thiouracil) FW 161.18 mp >300° Beil. 24,476 FT-IR 1(2):827A R&S 1(2):2577M RTECS# UW0495000	25g 100g	17.40 50.00
51,612-0	(S-(R:R'))-(+)-2-Amino-N-(2-hydroxy-1-methyl-2-phenylethyl)-N-methyl- acetamide, 97% [170115-96-5] $H_2NCH_2CON(CH_3)CH(CH_3)CH(OH)C_6H_5$ FW 222.29 mp 76-80° [α] _D ¹⁸ +101° (c=2, CH ₃ OH) MOISTURE-SENSITIVE	1g 5g	22.20 73.80
	2-Amino-2-(hydroxymethyl)-1,3-propanediol, see Tris(hydroxymethyl)- aminomethane		
A5,800-3	2-Amino-4-hydroxy-6-methylpyrimidine, 98% [3977-29-5] (2-amino-6-methyl- 4-pyrimidinol) FW 125.13 mp >300° Beil. 24,343 FT-IR 1(2):828B R&S 1(2):2579E	25g 100g	12.90 41.80
37,647-7	4-Amino-5-hydroxy-2,7-naphthalenedisulfonic acid, disodium salt, tech., 65% [3963-80-2] $H_2NC_{10}H_4(OH)(SO_3Na)_2$ FW 363.28 Merck Index 13,6305 FT-NMR 1(2):1604A R&S 1(2):2215K RTECS# QJ6190000 IRRITANT Contains 10-20% Na ₂ SO ₄	100g 500g	25.80 62.00
52,876-5	4-Amino-5-hydroxy-2,7-naphthalenedisulfonic acid, monosodium salt hydrate, 85% [312693-54-2] (H acid) $H_2NC_{10}H_4(OH)(SO_3H)SO_3Na \cdot xH_2O$ FW 341.30 mp >300° IRRITANT	100g 500g	40.40 122.50
39,896-9	4-Amino-3-hydroxy-1-naphthalenesulfonic acid, 98+%, A.C.S. reagent [116-63-2] $H_2NC_{10}H_5(OH)SO_3H$ FW 239.25 mp 290°(dec.) Beil. 14,846 Merck Index 13,453 FT-IR 1(2):500A R&S 1(2):2201L RTECS# QK1292000 CORROSIVE Assay ≥98.0% SO ₄ ²⁻ ≤0.2% Solubility in Na ₂ CO ₃ to pass Sens. to PO ₄ ³⁻ to pass Ignition residue ≤0.1%	5g 25g	11.80 36.20
10,975-4	4-Amino-3-hydroxy-1-naphthalenesulfonic acid, 95% [116-63-2] $H_2NC_{10}H_5(OH)SO_3H$	25g 100g	21.90 74.00
21,686-0	4-Amino-5-hydroxy-1-naphthalenesulfonic acid, tech., 90% [83-64-7] $H_2NC_{10}H_5(OH)SO_3H$ FW 239.25 Beil. 14,835 FT-IR 1(2):500B Safety 2,181A R&S 1(2):2201M RTECS# QK1295000 CORROSIVE	5g 100g	18.30 35.20
29,079-3	6-Amino-4-hydroxy-2-naphthalenesulfonic acid monohydrate, 90% [139123-65-2] $H_2NC_{10}H_5(OH)SO_3H \cdot H_2O$ FW 257.27 Beil. 14(3):2289 Safety 2,181B R&S 1(2):2201H CORROSIVE	5g 250g 1kg	19.10 35.30 111.90
28,254-5	7-Amino-4-hydroxy-2-naphthalenesulfonic acid monohydrate, 97% [139123-66-3] $H_2NC_{10}H_5(OH)SO_3H \cdot H_2O$ FW 257.27 Safety 2,181D R&S 1(2):2201J CORROSIVE	5g 100g 500g	19.10 34.50 109.70
51,522-1	3-Amino-4-hydroxy-5-nitrobenzenesulfonic acid, 97% [96-93-5] $H_2NC_6H_3(OH)(NO_2)SO_3H$ FW 234.19 mp 325°(dec.) Beil. 14,816 CORROSIVE	5g	49.10
	4-Amino-5-hydroxy-3-(4-nitrophenylazo)-2,7-naphthalenedisulfonic acid, disodium salt, see 20,890-6, Nitro Red page 1360		
	4-Amino-5-hydroxy-3-(4-nitrophenylazo)-6-(phenylazo)-2,7-naphthalene disulfonic acid, disodium salt, see 19,524-3, Naphthol Blue Black, page 1320		
	(R)-Amino-(4-hydroxyphenyl)acetic acid methyl ester hydrochloride, see 53,492-7; D-(-)-D-Hydroxyphenylglycine methyl ester hydrochloride page 1046		



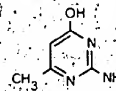
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A5,680-9

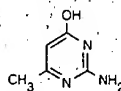


A5,740-6



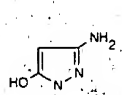
A5,800-3

	5g	40.80
(O)	25g	135.20
	1g	18.80
(O)(CH ₂ CO ₂ H)	5g	62.10
drochloride, atendine		
10:1-piperazine- B R&S 1(1),393M	1g	118.80
	5g	397.30
enol) FW 197.24 8410000	100mg	31.20
[65802-56-1]	25g	17.40
11.18 mp 300°	100g	50.00
(1)-N-methyl H ₂ C ₄ H ₈ IVE (ethyl)-	1g	22.20
	5g	73.80
ino-6-methyl- R&S 1(2),2579E	25g	12.90
	100g	41.80
-salt, tech., 65% 05	100g	25.80
	500g	62.00
m salt hydrate, FW 341.30	100g	40.40
	500g	122.50
reagent 346 Merck CORROSIVE	5g	11.80
	25g	36.20
2]	25g	21.90
	100g	74.00
3-64-7]	5g	18.30
2,181A	100g	35.20
90%	5g	19.10
39 Safety 2,181B	250g	35.30
	1kg	111.90
97%	5g	19.10
R&S 1(2),2201J	100g	34.50
	500g	109.70
3]	5g	49.10
CORROSIVE		
ifonic acid,		
phthalene black page 1320		
loride, see le page 1046		



A5,800-3

(S)-Amino-(4-hydroxyphenyl)acetic acid methyl ester hydrochloride, see 53,491-9; L-(+)-p-Hydroxyphenylglycine methyl ester hydrochloride page 1046			
48,260-9	N-[(2S,3R)-3-Amino-2-hydroxy-4-phenylbutyl]-L-leucine, 97%, [58970-76-6] (CH ₃) ₂ CHCH ₂ CH(NHCOCH(OH)CH(NH ₂)CH ₂ C ₆ H ₄ CO ₂ H FW 308.38 mp 245° (dec.) [α] _D ²⁰ -11° (c=1, 1N NaOH) Merck Index 13,9910	500mg	132.60
3-Amino-2-hydroxypropionic acid, see 28,633-8; DL-Isoserine page 1108			
4-Amino-4-(3-hydroxypropyl)-1,7-heptanediol, see 36,154-2; Bis-homotris page 227			
33,144-9	3-Amino-5-hydroxypyrazole, 98% [6126-22-3] FW 99.09 mp 214° (dec.) Beil 25(4),3516 FT-NMR 1(3),72A R&S 1(2),2347G CORROSIVE	5g	19.20
		25g	63.40
12,251-3	2-Amino-3-hydroxypyridine, 98% [16867-03-1] (2-amino-3-pyridinol) FW 110.12 mp 172-174° Beil 22(2),408 FT-IR 1(2),770A Safety 2,182A R&S 1(2),2527G IRRITANT	5g	12.10
		25g	39.20
		100g	143.80
		1kg	580.90
		5kg	1,828.30
08128	2-Amino-8-hydroxyquinoline [70125-16-5] FW 160.18 mp 157-160° Fluka	1g	43.35
		5g	165.25
30,552-9	5-Amino-8-hydroxyquinoline dinitrochloride, 95% [21302-43-2] (5-amino-8-quinolindol)dinitrochloride) FW 233.10 mp 279° (dec.) Beil 22(4),5866 R&S 1(2),2621F IRRITANT	1g	24.40
		5g	80.50
54,603-8	2-Amino-4-hydroxy-6-trifluoromethylpyrimidine, 97% [1513-69-5] FW 181.10 mp 289.5-293.5° IRRITANT	1g	98.40
19,791-2	2-Aminoimidazole sulfate, 98% [1450-93-7] FW 132.13 mp 270° (dec.) Beil 24(1),188 FT-IR 1(2),614D R&S 1(2),2353M IRRITANT	2.5g	65.70
		10g	171.50
55,241-0	5-Aminoimidazole-4-carboxamide, 95% [360-97-4] FW 126.12 mp 164-170° RTECS# N1390000 IRRITANT	25g	146.20
16,496-8	4-Amino-5-imidazolecarboxamide hydrochloride, 98% [72-40-2] FW 162.58 mp 250-252° (dec.) Beil 25(2),221 FT-IR 1(2),625A Safety 2,182B R&S 1(2),2365D RTECS# N13911000	1g	17.00
	Reactant for synthesis of heterocyclic compounds such as purines ^{1,2} and pyrimidines ³ . (1) J. Heterocycl. Chem. 1993, 30, 593. (2) J. Org. Chem. 1991, 56, 2139. (3) Heterocycles 1992, 34, 1133.	5g	68.80
		25g	204.70
37,674-4	2-Amino-4,5-imidazoledicarbonitrile, 97% [40953-34-2] FW 133.11 mp 270° (dec.) FT-NMR 1(3),88A R&S 1(2),2365I CORROSIVE	10g	43.40
		50g	143.70
Aminoiminomethanesulfonic acid, see F1,600-1; Formamidinesulfonic acid page 933			
N-(Aminoiminomethyl)-4-morpholinecarboximidamide hydrochloride, see 27,861-0; Moroxydine hydrochloride page 1311			
44,534-7	(R)-(-)-1-Aminoindan, 97% [10277-74-4] [(R)-(-)-1-indanamine] FW 133.19 mp 15° bp 96-97°/8mm n _D ²⁰ 1.5610 d 1.038 Fp 202°F(94°C) [α] _D ²⁰ -16.5° (c=1.5, CH ₃ OH) Beil 12(2),651 IRRITANT	250mg	39.80
		1g	105.30
A5,950-6	1-Aminoindan, 98% [34698-41-4] (1-indanamine) FW 133.19 mp 1.5° bp 96-97°/8mm n _D ²⁰ 1.5610 d 1.038 Fp 202°F(94°C) Beil 12,1191, FT-NMR 1(2),593A FT-IR 1(1),1281A Safety 2,182C R&S 1(1),1479I IRRITANT	1g	24.80
44,535-5	(S)-(+)-1-Aminoindan, 97% [61341-86-4] [(S)-(+)-1-indanamine] FW 133.19 bp 96-97°/8mm n _D ²⁰ 1.5620 d 1.038 Fp 202°F(94°C) [α] _D ²⁰ +16.5° (c=1.5, CH ₃ OH) Beil 12(2),651 IRRITANT	250mg	39.80
		1g	105.30
A5,952-2	2-Aminoindan hydrochloride, 98% [2338-18-3] (2-indanamine hydrochloride) FW 169.66 mp 246-247° Beil 12,1196 FT-NMR 1(2),593C FT-IR 1(1),1281C Safety 2,182D R&S 1(1),1479K RTECS# NK4050000 TOXIC	250mg	13.30
		1g	38.20
		5g	152.50
		25g	443.00
16,210-8	4-Aminoindan, 97% [32202-61-2] (4-indanamine) FW 133.19 bp 90-92° n _D ²⁰ 1.5920 d 0.736 Fp >230°F(110°C) IRRITANT	1g	5.80
		5g	18.90



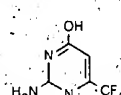
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12,251-3



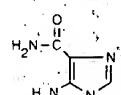
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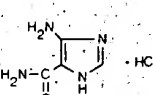
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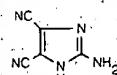
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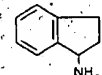
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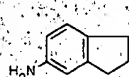
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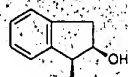
A5,950-6

■ Aminoindan ■

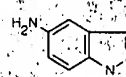
			US \$
13,087-7	5-Aminoindan, 95% [24425-40-9] (5-indanamine) FW 133.19 mp 34-36° bp 247-249°/745mm Fp >230°F(110°C) Beil. 12(1),511 FT-NMR 1(2),495B FT-IR 1(1):1215D Safety 2,183A R&S 1(1),1405E IRRITANT	1g 10g 50g	9.50 42.20 174.00
44,084-1	(1R,2S)-(+)-cis-1-Amino-2-indanol, 99% [136030-00-7] FW 149.19 mp 118-121° [α] _D ²⁵ +63° (c=0.2, CHCl ₃) 99% ee/GLC	1g 5g	26.90 88.90
44,083-3	(1S,2R)-(-)-cis-1-Amino-2-indanol, 99% [126456-43-7] FW 149.19 mp 118-121° [α] _D ²⁵ -61° (c=0.5, CHCl ₃) 99% ee/GLC	1g 5g	30.00 99.40
A5,955-7	5-Aminoindazole, 97% [19335-11-6] FW 133.15 mp 175-178° Beil. 25(2),308 ★ FT-NMR 1(3),176A FT-IR 1(2),686C Safety 2,183B R&S 1(2),2437C RTECS# NK7711000 TOXIC IRRITANT	5g 25g	33.50 105.30
A5,956-5	6-Aminoindazole, 98% [6967-12-0] FW 133.15 mp 204-206°(dec.) Beil. 25,317 ★ FT-NMR 1(3),176B FT-IR 1(2),686D Safety 2,183C R&S 1(2),2437D RTECS# NK7712000 IRRITANT	5g 25g	17.50 57.40
52,502-2	4-Aminoindole, 97% [5192-23-4] FW 132.17 mp 106-109° IRRITANT	500mg	86.30
A5,965-4	5-Aminoindole, 97% [5192-03-0] FW 132.17 mp 131-133°(dec.) Beil. 22(4),4296 ★ FT-IR 1(2),661D R&S 1(2),2407A	250mg 1g 5g	18.90 52.30 177.50
30,720-3	5-Aminoindole hydrochloride, 98% [65795-92-8] FW 168.63 mp 255-257° FT-NMR 1(3),133B R&S 1(2),2407B	250mg 1g	27.70 77.20
A5,960-3	2-Amino-5-iodobenzoic acid, 97% [5326-47-6] (5-iodoanthranilic acid) ★ H ₂ NC ₆ H ₃ (I)CO ₂ H FW 263.03 mp 219-221°(dec.) Beil. 14,373 FT-NMR 1(2),1121A FT-IR 1(2),219A Safety 2,184B R&S 1(2),1807N RTECS# DG2802000 IRRITANT	5g 25g	19.30 64.60
51,639-2	2-Amino-5-iodopyridine, 98% [20511-12-0] FW 220.01 mp 128-131° Beil. 22(2),334 ★ IRRITANT	1g	28.40
85,099-3	2-Aminoisobutyric acid, 98% [62-57-7] (2-methylalanine) (CH ₃) ₂ C(NH ₂)CO ₂ H ★ FW 103.12 mp >300° Beil. 4,414 Merck Index 13,444 FT-NMR 1(1),870B FT-IR 1(1),571C R&S 1(1),653N RTECS# AY7000000	25g 100g	18.50 55.90
21,779-4	DL-3-Aminoisobutyric acid, 98% [10569-72-9] H ₂ NCH ₂ CH(CH ₃)CO ₂ H FW 103.12 mp 179-182° Beil. 4(3),1330 FT-NMR 1(1),871A FT-IR 1(1),572B R&S 1(1),655A	1g 5g	27.40 108.70
41,592-8	3-Amino-1H-indole hydrochloride [76644-74-1] FW 168.63 mp 130°(dec.) IRRITANT, HYGROSCOPIC	1g 5g	21.70 70.70
18,627-9	5-Aminoisophthalic acid, 94% [99-31-0] H ₂ NC ₆ H ₃ -1,3-(CO ₂ H) ₂ FW 181.15 ★ mp >300° Beil. 14(1),636 FT-IR 1(2),228A R&S 1(2),1817D IRRITANT	25g 100g	26.00 71.90
40,230-3	2-Amino-7-isopropyl-5-oxo-5H-[1]benzopyrano[2,3-b]pyridine-3-carbonitrile, 99% [68302-12-5] FW 279.30 mp >300° IRRITANT	1g	29.80
45,064-2	3-Amino-2-isopropyl-4(3H)-quinazolinone, 99% [70589-51-4] FW 203.25 mp 104-106° IRRITANT	1g 5g	36.80 122.00
17,859-4	1-Aminoisoquinoline, 99% [1532-84-9] FW 144.18 mp 122-124° Beil. 22(1),640 ★ FT-NMR 1(3),458C R&S 1(2),2641H IRRITANT	1g 5g	39.70 154.30
13,610-7	5-Aminoisoquinoline, 99% [1125-60-6] FW 144.18 mp 125-128° Beil. 22,452 ★ FT-NMR 1(3),459A FT-IR 1(2),879A R&S 1(2),2641H IRRITANT	1g 5g	27.90 89.50
42,421-8	3-Aminoisoxazole, 95% [1750-42-1] FW 84.08 bp 75-76° n _D ²⁰ 1.5110 d 1.138 Fp >230°F(110°C) Beil. 27(4),4572 IRRITANT	5mL 25mL	29.00 116.60
	4-Amino-3-isoxazolidinone, see Cycloserine		
A5,990-5	5-Aminolevulinic acid hydrochloride, 99% [5451-09-2] H ₂ NCH ₂ COCH ₂ CH ₂ CO ₂ H·HCl FW 167.59 mp 156°(dec.) Merck Index 13,445 ★ FT-IR 1(1),582C R&S 1(1),663A RTECS# OI1640000 IRRITANT Substrate for 5-aminolevulinic acid dehydratase. Anal. Biochem. 1974, 58, 500. Clin. Chem. 1974, 20, 753, 1287.	100mg 500mg	17.50 55.50



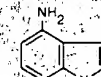
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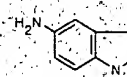
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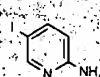
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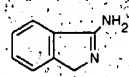
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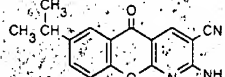
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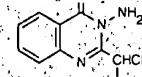
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41,592-8



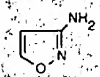
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45,064-2



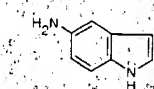
17,859-4



42,421-8

mp 34-36°	1g	9.50	6-Amino-2,4-lutidine, see A5,180-7, 2-Amino-4,6-dimethylpyridine page 81		
1,3,5-AB	10g	42.20	22,114-7 Aminomalonalonitrile <i>p</i> -toluenesulfonate, 98% [5098-11-6]	5g	65.20
	50g	174.00	H ₂ NCH(CN)·C ₁₀ H ₇ SO ₃ H FW 253.28 mp 174°(dec.) FT-NMR 1(2),1598A	25g	216.60
19 mp 118-121°	1g	26.90	FT-IR 1(2),502B Safety 2,184C R&S 1(2),2207D		
	5g	88.90	53,606-7 5-Amino-2-mercaptobenzimidazole, 96% [2818-66-8] FW 165.21 mp 240-244°	25g	53.80
19 mp 118-121°	1g	30.00	★ IRRITANT	100g	149.30
	5g	99.40	45,437-0 6-Amino-2-mercaptobenzothiazole, 97% [7442-07-1] (6-amino-2-benzothiazole-2-thiol) FW 182.27 mp 267-272° RTECS# D16488000 IRRITANT	1g	31.90
21,25(2),30B	5g	33.50	★	5g	106.70
	25g	105.30	2-Amino-4-mercaptobutyric acid, see 19,314-3, DL-Homocysteine page 1004		
1, Beil 25,316	5g	17.50	85,841-2 (-)-2-Amino-6-mercaptapurine riboside hydrate, 98% [40668-18-6] (thio-	250mg	24.40
	25g	57.40	guanosine) FW 299.31 mp 230-231°(dec.) [α] _D ²⁰ -69.9° (c=1.3, 0.1N NaOH) Merck	1g	78.60
ANI	500mg	86.30	Index 13,9267 FT-IR 1(2),718C Safety 2,184D R&S 1(2),2475D RTECS# UP0200000		
Beil 22(4),4296	250mg	18.90	A6,020-2 4-Amino-6-mercaptopyrazolo[3,4-d]pyrimidine, 95% [23771-52-0] (4-amino-	1g	67.20
	1g	52.30	pyrazolo[3,4-d]pyrimidine-6-thiol) FW 167.19 mp >300° FT-IR 1(2),716A		
	5g	177.50	23,362-5 4-Amino-2-mercaptopyrimidine, 97% [333-19-3] (thiocytosine) FW 127.17	1g	29.40
p 255-257°	250mg	27.70	mp 184°(dec.) Beil 1,583 FT-IR 1(1),889D Safety 2,185A R&S 1(1),1059E	5g	114.30
ilic acid)	1g	77.20	RTECS# UW0520000		
4R 1(2),1121A	5g	19.30	6-Amino-2-mercapto-4-pyrimidinol monohydrate, see A5,740-6, 4-Amino-6-		
10, IRRITANT	25g	64.60	hydroxy-2-mercaptopyrimidine monohydrate page 88		
31° Beil 22(2),334	1g	28.40	14,026-0 3-Amino-5-mercapto-1,2,4-triazole, 95% [16691-43-3] (3-amino-1,2,4-triazole-	10g	33.40
			5-thiol) FW 116.15 mp >300° FT-IR 1(2),628B R&S 1(2),2367L RTECS# XZ5357500	50g	131.00
			12,744-2 Aminomethanesulfonic acid, 97% [13881-91-9] H ₂ NCH ₂ SO ₃ H FW 111.12	10g	10.80
			mp 184°(dec.) Beil 1,583 FT-IR 1(1),889D Safety 2,185A R&S 1(1),1059E	50g	29.80
			★ CORROSIVE		
			Aminomethanetrimethanol, see Tris(hydroxymethyl)aminomethane		
			24,666-2 3'-Amino-4'-methoxyacetanilide, tech., 90% [6375-47-9]	25g	32.20
			★ CH ₃ CONHC ₆ H ₄ (NH ₂)OCH ₃ FW 180.21 Beil 13(3),1341 FT-IR 1(2),359B Safety 2,185B		
			R&S 1(2),1999D IRRITANT		
			41,594-4 2-Amino-4'-methoxyacetophenone hydrochloride, 90%, tech., [3883-94-1]	1g	35.40
			★ CH ₃ OC ₆ H ₄ COCH ₂ NH ₂ ·HCl FW 201.66 mp 190-193° Beil 14(3),549 IRRITANT	5g	116.80
			15,703-1 3-Amino-4-methoxybenzenesulfonic acid, 98% [98-42-0] (4-methoxymetani-	25g	132.60
			★ acid) H ₂ NC ₆ H ₃ (OCH ₃)SO ₃ H FW 203.22 mp 305°(dec.) Beil 14,814 CORROSIVE		
			34,010-3 2-Amino-3-methoxybenzoic acid, 98% [3177-80-8] (3-methoxyanthranilic acid)	1g	30.90
			★ H ₂ NC ₆ H ₃ (OCH ₃)CO ₂ H FW 167.16 mp 169-170° Beil 14(1),654 FT-NMR 1(2),1123C	5g	101.10
			R&S 1(2),1809H IRRITANT	25g	335.70
			18,606-6 3-Amino-4-methoxybenzoic acid, 98% [2840-26-8] H ₂ NC ₆ H ₃ (OCH ₃)CO ₂ H	10g	43.70
			★ FW 167.16 mp 208-210° Beil 14(1),657 FT-NMR 1(2),1124A FT-IR 1(2),220B		
			Safety 2,185C R&S 1(2),1809I RTECS# DG2872000 IRRITANT		
			34,702-7 4-Amino-3-methoxybenzoic acid, 98% [2486-69-3] H ₂ NC ₆ H ₃ (OCH ₃)CO ₂ H	1g	41.60
			★ FW 167.16 mp 186-188° Beil 14(1),654 FT-NMR 1(2),1124B R&S 1(2),1809J	5g	162.70
			IRRITANT		
			13,821-5 2-Amino-4-methoxybenzothiazole, 97% [5464-79-9] FW 180.23 mp 153-155°	1g	31.00
			★ FT-NMR 1(3),208A FT-IR 1(2),703D R&S 1(2),2457D RTECS# DL2000000	5g	113.80
				25g	359.80
			16,259-0 2-Amino-6-methoxybenzothiazole, 98% [1747-60-0] FW 180.23 mp 165-167°	50g	59.10
			★ Beil 27(2),334 FT-NMR 1(3),208B FT-IR 1(2),704A Safety 2,185D R&S 1(2),2457E	250g	213.20
			RTECS# DL2100000 TOXIC		
			A6,080-6 2-Amino-6-methoxybenzothiazole, tech., [1747-60-0]	25g	15.10
			★	100g	41.90

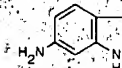
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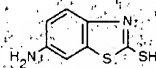
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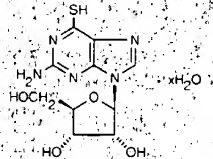
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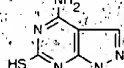
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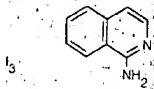
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A6,020-2



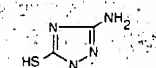
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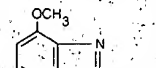
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42,421-8



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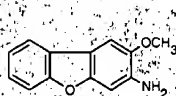
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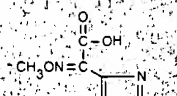
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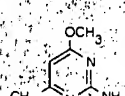
			US \$	
26,455-5	3-Amino-2-methoxydibenzofuran, 97% [5834-17-3] FW 213.24 mp 92-94°	25g	42.00	
	Beil. 18(4), 7352 FT-NMR 1(3), 172B. Safety 2, 186A R&S 1(2), 2431O	100g	139.00	
	RTECS# HP4900000 IRRITANT			
28,014-3	2-Amino-4-(methoxymethyl)-4-thiazoleacetic acid, 97%, predominantly syn	5g	38.70	
	[65872-41-5] FW 201.21 mp 192°(dec.) Safety 2, 186B R&S 1(2), 2391G	25g	118.80	
	MOISTURE-SENSITIVE IRRITANT			
	4'-Amino-5'-methoxy-2'-methylbenzanilide, see 20,158-8, Fast Violet B			
	page 894			
37,946-8	2-Amino-4-methoxy-6-methylpyrimidine, 98% [7749-47-5] FW 139.16	5g	32.40	
	mp 156-158° Beil. 25(4), 3385 FT-NMR 1(3), 381A R&S 1(2), 2575J IRRITANT	25g	105.60	
28,158-1	(R)-(+)-1-Amino-2-(methoxymethyl)pyrrolidine, 96% [72748-99-3] (RAMP)	250mg	37.60	
	FW 130.19 bp 42°/1.8mm n _D 1.4650 d 0.970 Fp 162°F(72°C) [α] _D ¹⁸ +79° (neat)	1g	103.70	
	Fieser 12.30 17.15 Safety 2, 186D R&S 1(1), 379K IRRITANT			
	Reagent for asymmetric synthesis			
	99% ee/GLC			
28,157-3	(S)-(-)-1-Amino-2-(methoxymethyl)pyrrolidine, 95% [59983-39-0] (SAMP)	1g	104.80	
	FW 130.19 bp 42°/1.8mm n _D 1.4640 d 0.970 Fp 162°F(72°C) [α] _D ¹⁸ -79° (neat)			
	Fieser 12.30 14.22 16.12 17.15 Safety 2, 186C R&S 1(1), 379L IRRITANT			
	Undergoes condensation reactions with carbonyl compounds to produce hydrazones			
	which are useful synthons in highly enantioselective syntheses. Tetrahedron 1993, 49			
	1821. Synthesis 1993, 725. Synlett 1992, 897			
	97% ee/GLC			
52,299-6	2-Amino-4-methoxy-6-methyl-1,3,5-triazine, 97% [1668-54-8] FW 140.15	10g	50.50	
	mp 258-261° IRRITANT			
26,211-0	5-Amino-2-methoxyphenol, 98% [1687-53-2] (3-hydroxy-4-methoxyaniline)	1g	19.30	
	H ₂ NC ₆ H ₃ (OCH ₃)OH FW 139.16 mp 130-132° Beil. 13(1), 307 FT-NMR 1(2), 514A	10g	106.40	
	Safety 2, 187A R&S 1(1), 1415I IRRITANT			
54,917-7	5-Amino-3-(4-methoxyphenyl)-1-phenylpyrazole, 97% [19652-13-2] FW 265.31	1g	30.30	
	mp 192-196° IRRITANT	5g	101.00	
31,652-0	(1S,2S)-(+)-2-Amino-3-methoxy-1-phenyl-1-propanol, 98% [51594-34-4]	250mg	32.50	
	CH ₃ OCH ₂ CH(NH ₂)CH(C ₆ H ₅)OH FW 181.24 mp 51-53° Fp >230°F(110°C)	1g	89.50	
	[α] _D ¹⁸ +25° (c=10.6, CHCl ₃) Fieser 12,381 13.17 FT-NMR 1(2), 589B R&S 1(1), 1477E			
	IRRITANT			
53,537-0	3-Amino-5-(4-methoxyphenyl)pyrazole, 97% [179541-95-8] FW 189.22	1g	34.70	
	mp 140-144° IRRITANT			
A6,100-4	2-Amino-1-methoxypropane, 95% [37143-54-7] (2-methoxyisopropylamine)	5g	23.40	
	CH ₃ OCH ₂ CH(CH ₃)NH ₂ FW 89.14 bp 92.5-93.5°/743mm n _D 1.4060 d 0.845	100g	45.40	
	Fp 48°F(8°C) [α] _D ²⁵ 0° (c=1, CHCl ₃) Beil. 4(4), 1615 FT-NMR 1(1), 526A Safety 2, 187B			
	R&S 1(1), 355E FLAMMABLE LIQUID - CORROSIVE			
	2-Amino-6-methoxypurine, see 36,305-7, O-Methylguanine page 1245			
A6,120-9	5-Amino-2-methoxypyridine, 95% [6628-77-9] FW 124.14 mp 29-31°	5g	32.70	
	bp 85-90°/11mm n _D 1.5750 Fp >230°F(110°C) Beil. 22(2), 408 FT-NMR 1(3), 300B	25g	109.20	
	FT-IR 1(2), 772D Safety 2, 187C R&S 1(2), 2529F RTECS# US1836000 IRRITANT			
51,324-5	4-Amino-6-methoxypyrimidine, 98% [696-45-7] FW 125.13 mp 157-160°	5g	49.80	
	Beil. 25(4), 3359 IRRITANT			
52,809-9	8-Amino-6-methoxyquinoline hydrobromide, 97% [312693-53-1] FW 255.11	1g	66.10	
	mp 238°(dec.) IRRITANT	10g	366.60	



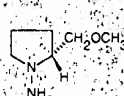
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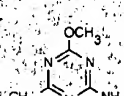
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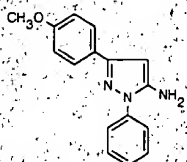
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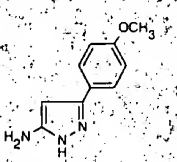
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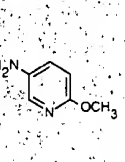
52,299-6



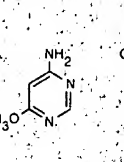
54,917-7



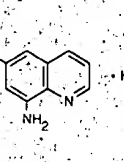
53,537-0



A6,120-9



51,324-5



52,809-9

-94°	25g	US \$ 42.00
.....	100g	139.00
ly. syn	5g	38.70
.....	25g	118.80

it B		
3	5g	32.40
VT	25g	105.60
AMP)	250mg	37.60
eat)	1g	103.70

AMP)	1g	104.80
eat)		

frazones
1993, 49,

1.15	10g	50.50
------------	-----	-------

aniline)	1g	19.30
1,514A	10g	106.40

FW 265.31	1g	30.30
.....	50g	101.00

4-4)	250mg	32.50
.....	1g	89.50

l(1), 1477E		
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2	1g	34.70
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amine)	5g	23.40
.....	100g	45.40

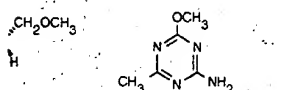
ety 2, 187B

45

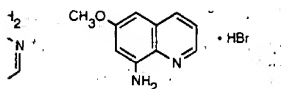
.....	5g	32.70
.....	25g	109.20

.....	5g	49.80
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V 255.11	1g	66.10
.....	10g	366.60



158-1 52,299-6



5 52,809-9

10-962-9591

1-(Aminomethyl)adamantane, see 18,037-8, 1-Adamantanemethylamine page 40

1-Amino-2-methylantraquinone, see 21,709-3, Disperse Orange 11 page 792

Aminomethylated polystyrene, see StratoSpheres™ PL-AMS resin, or ScavengePore® aminomethylated polystyrene

Aminomethylated styrene/divinylbenzene copolymer, see Poly(styrene-co-divinylbenzene), aminomethylated

54,825-1	3-Amino-4-methylbenzamide, 97% [19406-86-1] NH ₂ C ₆ H ₃ (CH ₃)CONH ₂ FW 150.18	5g	26.10
★	mp 127-131° IRRITANT	25g	86.80

A6,180-2	4-(Aminomethyl)benzenesulfonamide hydrochloride hydrate, 99% [138-37-4] (α-amino-p-toluenesulfonamide hydrochloride, homösulfonylamide hydrochloride) H ₂ NCH ₂ C ₆ H ₄ SO ₂ NH ₂ ·HCl·xH ₂ O FW 222.69 mp 261-263° Merck Index 13,5671 FT-NMR 1(2),1637B FT-IR 1(2),532C R&S 1(2),2243D RTECS# XT5425000	25g	28.70
		100g	100.20

Aminomethylbenzenesulfonic acid, see Aminotoluenesulfonic acid

16,563-8	2-(Aminomethyl)benzimidazole dihydrochloride hydrate, 98% [5993-91-9] FW 220.10 mp 264°(dec.) FT-NMR 1(2),184A FT-IR 1(2),692A R&S 1(2),2441G	5g	44.10
		25g	144.10

41,254-6	2-Amino-1-methylbenzimidazole, 95% [1622-57-7] FW 147.18 Beil. 25(5),10,408 RTECS# DD5423660 IRRITANT	1g	31.20
		5g	103.50

28,374-6	4-(Aminomethyl)benzoic acid, 97% [56-91-7] H ₂ NCH ₂ C ₆ H ₄ CO ₂ H FW 151.17 Beil. 14(3),1212 FT-NMR 1(2),1085A Safety 2,188D R&S 1(2),1793C IRRITANT	5g	21.50
		25g	71.30
		100g	197.10

A6,220-5	2-Amino-3-methylbenzoic acid, 99% [4389-45-1] (3-methylantranilic acid) H ₂ NC ₆ H ₃ (CH ₃)CO ₂ H FW 151.17 mp 176-178° Beil. 14,479 FT-NMR 1(2),1097B FT-IR 1(2),206A R&S 1(2),1797M IRRITANT	5g	41.60
		25g	137.10

41,944-3	2-Amino-5-methylbenzoic acid, 99% [2941-78-8] (5-methylantranilic acid) H ₂ NC ₆ H ₃ (CH ₃)CO ₂ H FW 151.17 mp 175°(dec.) Beil. 14,481 FT-IR 1(2),217A IRRITANT	1g	32.70
		5g	108.50
		25g	339.70

23,053-7	2-Amino-6-methylbenzoic acid, 99% [4389-50-8] (6-methylantranilic acid) H ₂ NC ₆ H ₃ (CH ₃)CO ₂ H FW 151.17 mp 127°(dec.) Beil. 14(1),598 FT-NMR 1(2),1097C FT-IR 1(2),206B Safety 2,188C R&S 1(2),1797N IRRITANT	5g	66.90
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33,418-9	3-Amino-2-methylbenzoic acid, 97% [52130-17-3] H ₂ NC ₆ H ₃ (CH ₃)CO ₂ H FW 151.17 mp 178-181° FT-NMR 1(2),1098A R&S 1(2),1797O IRRITANT	10g	29.10
		50g	105.70
		250g	321.70

A6,280-9	3-Amino-4-methylbenzoic acid, 99% [2458-12-0] H ₂ NC ₆ H ₃ (CH ₃)CO ₂ H FW 151.17 mp 167-169° Beil. 14,487 FT-NMR 1(2),1117C FT-IR 1(2),217B R&S 1(2),1807D IRRITANT	50g	64.00
		250g	213.20

A6,300-7	4-Amino-3-methylbenzoic acid, 98% [2486-70-6] H ₂ NC ₆ H ₃ (CH ₃)CO ₂ H FW 151.17 mp 169-171° Beil. 14,480 FT-NMR 1(2),1118A FT-IR 1(2),217C R&S 1(2),1807E IRRITANT	1g	31.10
		10g	201.00

21,428-0	2-Amino-4-methylbenzonitrile, 97% [26830-96-6] H ₂ NC ₆ H ₃ (CH ₃)CN FW 132.17 mp 92-95° Beil. 14,485 FT-IR 1(2),457C IRRITANT	1g	39.30
		5g	130.30

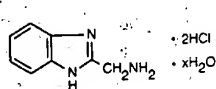
08419	2-Amino-6-methylbenzonitrile, pract., ≥95% (NT) [56043-01-7] FW 132.16 mp 127-132°	5g	59.40
Fluka		25g	234.90

54,958-4	5-Amino-2-methylbenzonitrile, 97% [50670-64-9] H ₂ NC ₆ H ₃ (CH ₃)CN FW 132.17 mp 88-91° IRRITANT	5g	38.30
		25g	130.20

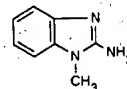
24,823-1	2-Amino-4-methylbenzophenone, 99% [4937-62-6] H ₂ NC ₆ H ₃ (CH ₃)COC ₆ H ₅ FW 211.26 mp 65-66° Beil. 14,107 FT-NMR 1(2),900C FT-IR 1(2),69B Safety 2,189B R&S 1(2),1675E IRRITANT	250mg	29.60
		1g	81.40

19,322-4	2-Amino-4-methylbenzothiazole, 97% [1477-42-5] FW 164.23 mp 137-139° Beil. 27,193 FT-NMR 1(3),206B FT-IR 1(2),702D Safety 2,189C R&S 1(2),2455J RTECS# DL2275000 TOXIC MOISTURE-SENSITIVE	5g	18.70
★		25g	45.50
		100g	147.70

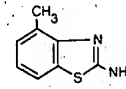
28,837-3	2-Amino-6-methylbenzothiazole, 99% [2536-91-6] FW 164.23 mp 140-142° Beil. 27(2),241 FT-NMR 1(3),206C Safety 2,189D R&S 1(2),2455K RTECS# DL2625000 IRRITANT	25g	24.30
★		100g	70.30



16,563-8



41,254-6

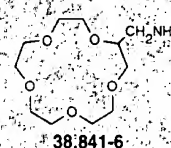
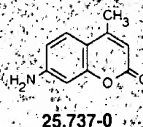


19,322-4

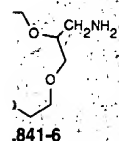
C of A's / MSDS's and spectra now available online at www.sigma-aldrich.com

Aminomethy

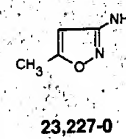
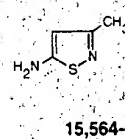
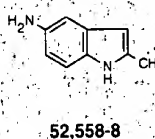
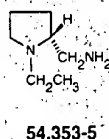
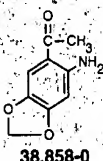
			US \$
A6,330-9	5-Amino-2-methylbenzothiazole dihydrochloride [32770-99-3] FW 237.15	5g	76.60
	mp 249-250° (dec.) Beil. 27(2), 427. FT-IR 1(2), 703A R&S 1(2), 2455L		
	RTECS# DL2650000		
	4-Aminomethylbenzoi 2-chlorotriyl resin, see 54,379-9, 2-Chlorotriyl 4-amino-		
	methylbenzoic acid, polymer-bound, page 478		
	(Aminomethyl)benzyl alcohol, see A7,240-5, 2-Amino-1-phenylethanol		
	page 103		
33,419-7	2-Amino-3-methylbenzyl alcohol, 97% [57772-50-6] H ₂ NC ₆ H ₃ (CH ₃)CH ₂ OH	5g	88.40
	FW 137.18 mp 67-69° FT-NMR 1(2), 493C R&S 1(1), 1403L IRRITANT	25g	294.50
33,298-4	2-Amino-5-methylbenzyl alcohol, 98% [34897-84-2] H ₂ NC ₆ H ₃ (CH ₃)CH ₂ OH	1g	30.70
	FW 137.18 mp 123-126° FT-NMR 1(2), 516C R&S 1(1), 1417D IRRITANT	5g	101.60
	3-Amino-4-methylbenzyl alcohol, see H2,620-0, 3-(1-Hydroxyethyl)aniline		
	page 1026		
33,420-0	3-Amino-2-methylbenzyl alcohol, 97% [83647-42-1] H ₂ NC ₆ H ₃ (CH ₃)CH ₂ OH	1g	17.60
	FW 137.18 mp 106-108° FT-NMR 1(2), 494A R&S 1(1), 1405A IRRITANT	10g	69.80
		50g	231.70
33,532-0	3-Amino-4-methylbenzyl alcohol, 97% [81863-45-8] H ₂ NC ₆ H ₃ (CH ₃)CH ₂ OH	5g	32.00
	FW 137.18 mp 107-109° Beil. 13,635 FT-NMR 1(2), 517A R&S 1(1), 1417E IRRITANT	25g	85.60
	2-Amino-3-methylbutane, see 19,018-7, 1,2-Dimethylpropylamine, page 753		
28,448-3	(R)-(-)-2-Amino-3-methyl-1-butanol, 98% [4276-09-9] (D-valinol)	1g	40.20
	(CH ₃) ₂ CHCH(NH ₂)CH ₂ OH. FW 103.17 mp 35-36° Fp 194°F(90°C)	5g	158.70
	[α] _D ²⁵ -16° (c=10, C ₂ H ₅ OH) Beil. 4(3), 805 Safety 2, 190C R&S 1(1), 363J IRRITANT		
18,483-7	2-Amino-3-methyl-1-butanol, 97% [473-75-6] (DL-valinol)	1g	17.80
	(CH ₃) ₂ CHCH(NH ₂)CH ₂ OH FW 103.17 bp 75-77°/8mm n _D ²⁰ 1.4540 d 0.936	5g	52.90
	Fp 194°F(90°C) [α] _D ²² 0° (neat) Beil. 4(2), 747 FT-IR 1(1), 336C Safety 2, 190A		
	R&S 1(1), 363L IRRITANT		
18,670-8	(S)-(+)-2-Amino-3-methyl-1-butanol, 96% [2026-48-4] (L-valinol)	1g	27.50
	(CH ₃) ₂ CHCH(NH ₂)CH ₂ OH FW 103.17 mp 30-32° bp 81°/8mm n _D ²⁰ 1.4540 d 0.926	5g	100.40
	Fp 196°F(91°C) [α] _D ²⁵ +10° (c=10, H ₂ O) Beil. 4(3), 805 Fieser 12,563 13,341, 16,380	25g	397.60
	FT-NMR 1(1), 540B FT-IR 1(1), 336D Safety 2, 190B R&S 1(1), 363K IRRITANT		
	Reacts with aldehydes and nitriles to form imines ¹ and oxazolines ² respectively, for		
	asymmetric synthesis. (1), J. Chem. Soc., Perkin Trans. 1 1989, 192. (2) Tetrahedron		
	Lett. 1993, 34, 2015, 3149-7793		
	97% ee/GLC		
	8-(4-Amino-1-methylbutylamino)-6-methoxyquinoline diphosphate, see		
	16,039-3, Primaquine diphosphate, page 1583		
	4-(Aminomethyl)catechol hydrobromide, see 85,878-1, 3,4-Dihydroxy-		
	benzylamine hydrobromide, page 681		
25,737-0	7-Amino-4-methylcoumarin, 99% [26093-31-2] (Coumarin 120) FW 175.19	100mg	29.00
	mp 223-226° λ _{max} 354nm Beil. 18,610 FT-NMR 1(2), 1318C FT-IR 1(2), 328D	500mg	84.20
	Safety 2: 190D R&S 1(2), 1945O IRRITANT, LIGHT-SENSITIVE		
	Fluorescent labeling reagent for trace determination of enzymes. For references, see		
	Aldrichimica Acta 1982, 15, 42		
	Suitable as laser dye		
23,773-6	7-Amino-4-methylcoumarin, 98% [26093-31-2]	250mg	29.10
		1g	90.10
38,841-6	2-(Aminomethyl)-15-crown-5, 95% [83585-56-2] FW 249.31 bp 115°/0.08mm	250mg	87.20
	n _D ²⁰ 1.4800 d 1.134 Fp >230°F(110°C) [α] _D ²⁶ 0° (neat) IRRITANT	1g	243.10
38,843-2	2-(Aminomethyl)-18-crown-6, 95% [83585-61-9] FW 293.36 bp 135°/0.08mm	250mg	128.70
	n _D ²⁰ 1.4790 d 1.126 Fp >230°F(110°C) [α] _D ²⁵ 0° (c=2, CHCl ₃) IRRITANT	1g	337.70
85,765-3	trans-4-(Aminomethyl)cyclohexanecarboxylic acid, 97% [1197-18-8]	10g	23.90
	(transamic acid) H ₂ NCH ₂ C ₆ H ₁₀ CO ₂ H FW 157.21 mp >300° Merck Index 13,9645	50g	79.90
	FT-IR 1(1), 543A Safety 2, 191A R&S 1(1), 595D RTECS# GU8400000 IRRITANT	250g	276.80
	Used as a lysine analogue to characterize binding sites in plasminogen. Biochemistry		
	1994, 33, 3599. ibid. 1993, 32, 10936		
19,141-8	1-Aminomethyl-1-cyclohexanol hydrochloride, 98% [19968-85-5]	1g	17.80
	H ₂ NCH ₂ C ₆ H ₁₀ OH·HCl FW 165.67 mp 217-219° FT-IR 1(1), 318A R&S 1(1), 365H	5g	58.80



FW 237.15	5g	US \$ 76.60
lorotyl, 4-amino-		
ylethanol		
CH ₃ CH ₂ OH	5g	88.40
T	25g	294.50
CH ₃ CH ₂ OH	1g	30.70
4NT	5g	101.60
thyl)aniline		
CH ₃ CH ₂ OH	1g	17.60
INT	10g	69.80
	50g	231.70
CH ₃ CH ₂ OH	5g	32.00
417E IRRITANT	25g	85.60
ie page 753		
)	1g	40.20
J IRRITANT	5g	158.70
	1g	17.80
1.936	5g	52.90
2.190A		
)	1g	27.50
540 d 0.926	5g	100.40
13.341 16.380	25g	397.60
IRITANT		
spectively, for		
2) Tetradron		
hate, see		
hydroxy-		
FW 175.19	100mg	29.00
2).328D	500mg	84.20
ferences, see		
	250mg	29.10
	1g	90.10
15%0.08mm	250mg	87.20
	1g	243.10
35%0.08mm	250mg	128.70
T	1g	337.70
18-8]	10g	23.90
k Index 13.9645	50g	79.90
IRRITANT	250g	276.80
l. Biochemistry		
-5]	1g	17.80
S 1(1).365H	5g	58.80

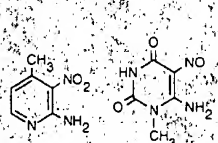


35,952-1	(Aminomethyl)cyclopropane, 97% [2516-47-4] (cyclopropanemethylamine) 1mL	US \$ 30.80
★	C ₃ H ₅ CH ₂ NH ₂ FW 71.12 bp 66°/758mm n _D ²⁰ 1.4340 d 0.820 Fp -23°F(-30°C) Beil. 12.4 FT-NMR 1(1).505B R&S 1(1).337B FLAMMABLE LIQUID CORROSIVE	
A6,380-5	(Aminomethyl)cyclopropane hydrochloride, 99% [7252-53-1] (cyclopropane-methylamine hydrochloride) C ₃ H ₅ CH ₂ NH ₂ ·HCl FW 107.58 mp 200° Beil. 12.4 FT-IR 1(1).312C Safety 2.191B R&S 1(1).337C IRRITANT HYGROSCOPIC	1g 24.30 10g 168.60
	α-(Aminomethyl)-3,4-dihydroxybenzyl alcohol, see Norepinephrine	
55,115-5	(R)-(+)-2-Amino-3-methyl-1,1-diphenylbutane, 98% FW 239.36 mp 73-76° 500mg	96.60
NEW	[α] _D ²⁵ +6.5° (c=1, CHCl ₃) IRRITANT A Product of Onyx Scientific, U.K.	
55,114-7	(S)-(-)-2-Amino-3-methyl-1,1-diphenylbutane, 98% FW 239.36 mp 73-76° 500mg	96.60
NEW	[α] _D ²⁵ -6.5° (c=1, CHCl ₃) IRRITANT A Product of Onyx Scientific, U.K.	
55,108-2	(R)-(+)-2-Amino-3-methyl-1,1-diphenyl-1-butanol, 98% [86695-06-9] 1g	103.60
NEW	(CH ₃) ₂ CHCH(NH ₂)C(C ₆ H ₅) ₂ (OH) FW 255.36 mp 95-99° [α] _D ²⁵ +132° (c=1, CHCl ₃) IRRITANT A Product of Onyx Scientific, U.K.	5g 344.70
55,101-5	(S)-(-)-2-Amino-3-methyl-1,1-diphenyl-1-butanol, 98% [78603-95-9] 1g	103.60
NEW	(CH ₃) ₂ CHCH(NH ₂)C(C ₆ H ₅) ₂ (OH) FW 255.36 mp 95-99° [α] _D ²⁵ -132° (c=1, CHCl ₃) IRRITANT A Product of Onyx Scientific, U.K.	5g 344.70
55,462-6	(S)-(-)-2-Amino-4-methyl-1,1-diphenylpentane, 97% 500mg	96.80
NEW	(CH ₃) ₂ CHCH ₂ CH(NH ₂)CH(C ₆ H ₅) ₂ FW 253.39 mp 48-52° Fp >230°F(110°C) [α] _D ²⁵ -32° (c=1, CHCl ₃) IRRITANT A Product of Onyx Scientific, U.K.	
55,455-3	(R)-(+)-2-Amino-4-methyl-1,1-diphenyl-1-pentanol, 98% [161832-74-2] 1g	103.70
NEW	(CH ₃) ₂ CHCH ₂ (NH ₂)C(C ₆ H ₅) ₂ OH FW 269.39 mp 144-147° [α] _D ²⁵ +100° (c=1, CHCl ₃) IRRITANT HYGROSCOPIC 99% ee/HPLC	5g 344.80
55,456-1	(S)-(-)-2-Amino-4-methyl-1,1-diphenyl-1-pentanol, 98% [78603-97-1] 1g	103.70
NEW	(CH ₃) ₂ CHCH ₂ CH(NH ₂)C(C ₆ H ₅) ₂ OH FW 269.39 mp 132-136° [α] _D ²⁵ -99° (c=1, CHCl ₃) IRRITANT HYGROSCOPIC	5g 344.80
38,858-0	6'-Amino-3',4'-(methylenedioxy)acetophenone, 98% [28657-75-2] FW 179.18 5g	12.00
✓	mp 170-174° Beil. 19(5).8.641 R&S 1(2).1657C IRRITANT	25g 39.50
54,353-5	(R)-(+)-2-Aminomethyl-1-ethylpyrrolidine, 96% [22795-97-7] FW 128.22 bp 50-52° 1g	41.50
NEW	n _D ²⁰ 1.4670 d 0.919 Fp 135°F(57°C) [α] _D ²⁵ -93° (c=1, CH ₃ OH) IRRITANT HYGROSCOPIC	5g 98.70
	2-Amino-6-methylheptane, see D16,129-2, 1,5-Dimethylhexylamine page 737	
29,620-1	6-Amino-2-methyl-2-heptanol hydrochloride, 99% [543-15-7] (heptaminol 25g	41.40
	hydrochloride) CH ₃ CH(NH ₂)(CH ₂) ₃ C(CH ₃) ₂ OH·HCl FW 181.71 mp 165-167° Merck Index 13.4676 Safety 2.191D R&S 1(1).365C RTECS# MJ3240000 IRRITANT HYGROSCOPIC	100g 105.00
	α-Aminomethyl-3-hydroxybenzyl alcohol hydrochloride, see 11,372-7, Norepinephrine hydrochloride page 1374	
	α-(Aminomethyl)-4-hydroxybenzyl alcohol hydrochloride, see 13,051-6, Octopamine hydrochloride page 1385	
	α-(Aminomethyl)-4-hydroxy-3-methoxybenzyl alcohol hydrochloride, see 28,688-5, Normetanephrine hydrochloride page 1374	
	4-(Aminomethyl)-5-hydroxy-6-methyl-3-pyridinemethanol dihydrochloride, see 28,709-1, Pyridoxamine dihydrochloride page 1610	
52,558-8	5-Amino-2-methylindole, 97% [7570-49-2] FW 146.19 mp 153-157° IRRITANT 5g	71.30
✓	15,564-0 5-Amino-3-methylisothiazole hydrochloride [52547-00-9] FW 150.63 mp 300° 5g	24.50
✓	FT-NMR 1(3).107C FT-IR 1(2).642C R&S 1(2).2385B RTECS# NX8150300 IRRITANT May contain ~10% NH ₄ Cl	25g 96.60
23,227-0	3-Amino-5-methylisoxazole, 98+%, [1072-67-9] FW 98.10 mp 61-63° 50g	29.30
✓	★ FT-NMR 1(3).99A FT-IR 1(2).636A Safety 2.192A R&S 1(2).2377H	250g 111.60

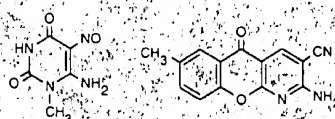


■ Aminomethy ■

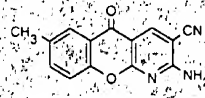
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30,427-1	5-Amino-3-methylisoxazole, 98% [14678-02-5] FW 98.10 mp 85-87° Beil. 27.157 ... FT-NMR 1(3).98A Safety 2.192B R&S 1(2).2377E IRRITANT	1g	33.10
		5g	103.80
29,008-4	2-Amino-4-methyl-3-nitropyridine, 98% [6635-86-5] FW 153.14 mp 139-141° Beil. 22(4).4180 FT-NMR 1(3).307C Safety 2.192C R&S 1(2).2533F IRRITANT	1g	16.40
		5g	54.40
29,009-2	2-Amino-4-methyl-5-nitropyridine, 98% [21901-40-6] FW 153.14 mp 223-225° Beil. 22(4).4180 FT-NMR 1(3).309A Safety 2.192D R&S 1(2).2533J IRRITANT	1g	29.60
56,696-9	6-Amino-1-methyl-5-nitrosouracil, 97% [6972-78-7] FW 170.13 mp >300° IRRITANT	25g	84.90
40,227-3	2-Amino-7-methyl-5-oxo-5H-[1]benzopyrano[2,3-b]pyridine-3-carbonitrile, 98% [68302-13-6] FW 251.25 mp >300° IRRITANT	1g	29.80
	2-Amino-3-methylpentanoic acid, see Isoleucine		
	2-Amino-3-methyl-1-pentanol, see 19,052-7, Isoleucinol page 1096		
	2-Amino-4-methyl-1-pentanol, see Leucinol		
	4-Amino-4-methyl-2-pentanone hydrogenoxalate, see 18,025-4, Diacetoneamine hydrogenoxalate page 561		
	Aminomethylphenol, see Aminocresol		
	2-(4-Amino-2-methylphenyl)-5-chloroisindole-1,3-dione, see 20,872-8, N-(4- Amino-2-methylphenyl)-4-chlorophthalimide page 96		
20,872-8	N-(4-Amino-2-methylphenyl)-4-chlorophthalimide, 99% [58230-69-6] FW 286.72 mp 208-210° FT-IR 1(2).410B R&S 1(2).2085B IRRITANT	10g	43.60
54,916-9	5-Amino-3-(4-methylphenyl)-1-phenylpyrazole, 97% [90012-40-1] FW 249.31 mp 173-176° IRRITANT	1g	37.90
		5g	126.30
54,100-1	5-Amino-3-methyl-1-phenylpyrazole, 97% [1131-18-6] FW 173.21 mp 114-117° RTECS# UO4990000 IRRITANT	5g	29.00
		25g	96.50
54,920-7	5-Amino-3-(4-methylphenyl)pyrazole, 97% FW 173.22 mp 148-151° IRRITANT	1g	34.70
		5g	115.70
34,631-4	2-[2-(Aminomethyl)phenylthio]benzyl alcohol, 95% [79467-22-4] H ₂ NCH ₂ C ₆ H ₄ SC ₆ H ₄ CH ₂ OH FW 245.34 mp 96-98° FT-NMR 1(2).597C R&S 1(1).1481M IRRITANT	5g	25.20
		25g	83.20
34,632-2	2-[2-(Aminomethyl)phenylthio]benzyl alcohol hydrochloride, 97% [62220-58-0] H ₂ NCH ₂ C ₆ H ₄ SC ₆ H ₄ CH ₂ OH · HCl FW 281.81 mp 197-199° R&S 1(1).1481N IRRITANT	5g	35.80
		25g	118.60
36,118-6	2-[2-(Aminomethyl)phenylthio]benzyl alcohol hydrochloride, tech., 90% [62220-58-0] H ₂ NCH ₂ C ₆ H ₄ SC ₆ H ₄ CH ₂ OH · HCl	5g	20.20
		25g	66.40
55,980-6	4-Amino-3-methyl-6-phenyl-1,2,4-triazin-5-one, 97% [41394-05-2] FW 202.21 mp 167-170° Merck Index 13.5949	25g	144.00
32,481-7	(Aminomethyl)phosphonic acid, 99% [1066-51-9] H ₂ NCH ₂ P(O)(OH) ₂ FW 111.04 Beil. 1(3).2596 FT-NMR 1(1).1475C R&S 1(1).1101I IRRITANT HYGROSCOPIC	250mg	28.00
		1g	111.30
24,771-5	4-Amino-N-methylphthalimide, 97% [2307-00-8] FW 176.18 mp 246-248° Beil. 22.535 FT-NMR 1(2).1464C FT-IR 1(2).410A R&S 1(2).2085A IRRITANT	250mg	20.60
		1g	57.40
25,568-8	1-Amino-4-methylpiperazine, 97% [6928-85-4] FW 115.18 bp 172-175° n _D ²⁰ 1.4850 d 0.957 Fp 145°F (62°C) Beil. 23(4).376 Safety 2.193B R&S 1(1).393K IRRITANT	25g	34.90
		100g	96.50
A6,513-1	1-Amino-4-methylpiperazine dihydrochloride monohydrate, 97% [40675-60-3] FW 206.12 mp 230° (dec.) FT-IR 1(1).374C Safety 2.193C R&S 1(1).393L RTECS# TK8062000 IRRITANT	1g	23.60
		10g	113.80



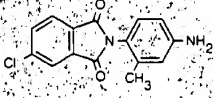
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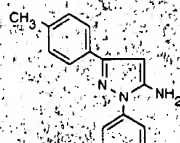
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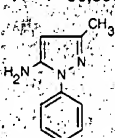
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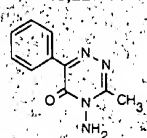
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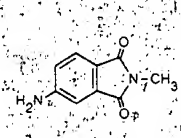
54,916-9



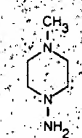
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55,980-6

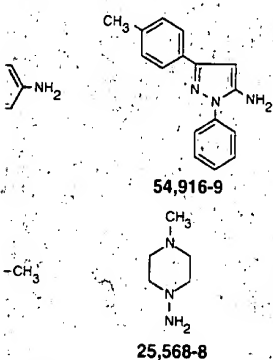


24,771-5

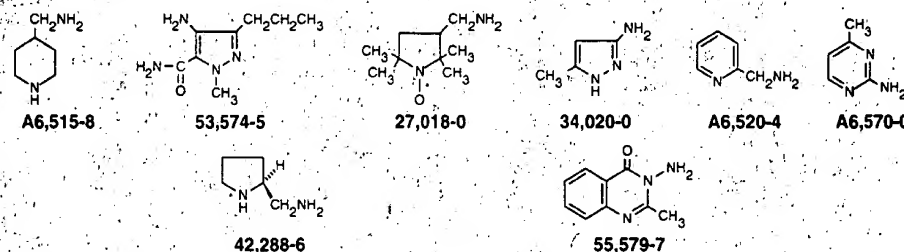


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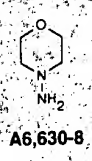
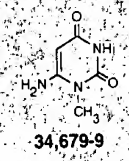
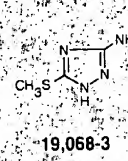
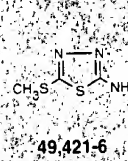
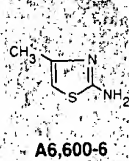
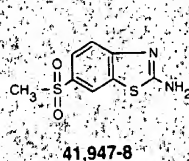
17° Beil. 27,157	1g	33.10
	5g	103.80
mp 139-141°	1g	16.40
IRITANT	5g	54.40
mp 223-225°	1g	29.60
IRITANT		
mp >300°	25g	84.90
rbonitrile, 98%	1g	29.80
096		
Diacetonamine		
0,872-8, N-(4-		
59-6] FW 286.72	10g	43.60
[] FW 249.31	1g	37.90
	5g	126.30
mp 114-117°	5g	29.00
	25g	96.50
1° IRITANT	1g	34.70
	5g	115.70
[5g	25.20
	25g	83.20
7%	5g	35.80
9°	25g	118.60
sch., 90%	5g	20.20
	25g	66.40
3] FW 202.21	25g	144.00
2H)₂ FW 111.04	250mg	28.00
IOSCOPIC	1g	111.30
46-248°	250mg	20.60
IRITANT	1g	57.40
175° n _D 1.4850	25g	34.90
IRITANT	100g	96.50
% [40675-60-3]	1g	23.60
3L	10g	113.80



A6,515-8	4-(Aminomethyl)piperidine, 96% [7144-05-0] FW 114.19 mp 25° bp 200°	5g	37.00
★	n _D 1.4900 Fp 174°F(78°C) FT-NMR 1(1),577A FT-IR 1(1),370B Safety 2,193D R&S 1(1),389G IRITANT		
	Aminomethyl, polymer-bound, see 51,477-2, Poly(styrene-co-divinylbenzene), aminomethylated page 1548		
A6,517-4	2-Amino-2-methyl-1,3-propanediol, 99% [115-69-5] HOCH ₂ C(CH ₃)(NH ₂)CH ₂ OH	5g	11.20
★	FW 105.14 mp 108-110° bp 151°/10mm Beil. 4,303 Fieser 2,24 Merck Index 13,447 FT-IR 1(1),347B Safety 2,194A R&S 1(1),371B RTECS# TY2975000 CORROSIVE	100g	38.50
		250g	67.80
A6,518-2	2-Amino-2-methyl-1-propanol, 95% [124-68-5] (CH ₃) ₂ C(NH ₂)CH ₂ OH FW 89.14	25mL	9.30
★	mp 24-28°, bp 165° n _D 1.4450 d 0.934 Fp 153°F(67°C) Beil. 4(3),783 Fieser 1,37 3,14 6,20 Merck Index 13,448 FT-NMR 1(1),539B Safety 2,194B R&S 1(1),363G	500mL	23.80
	RTECS# UA5950000 IRITANT	1L	31.50
	Used to derivatize carboxylic acids for GC analysis ¹ and to synthesize 2-oxazolines ² for further transformations. ³ (1) J. Chromatogr., A 1994, 673, 101. (2) Tetrahedron 1993, 49, 9353. (3) J. Am. Chem. Soc. 1992, 114, 1010.	12x1L	236.70
	May contain up to 5% monomethyl aminomethyl propanol		
47,186-0	(1R)-(+)-(1-Amino-2-methylpropyl)phosphonic acid, 98% [66254-56-6]	250mg	95.60
	(CH ₃) ₂ CHCH(NH ₂)P(O)(OH) ₂ FW 153.12 mp 272-277° [α] _D +1.0° (c=1, 1N NaOH) IRITANT		
47,187-9	(1S)-(-)-(1-Amino-2-methylpropyl)phosphonic acid, 98% [66254-55-5]	250mg	96.10
	(CH ₃) ₂ CHCH(NH ₂)P(O)(OH) ₂ FW 153.12 mp 272-277° [α] _D -1.0° (c=1, 1N NaOH) IRITANT		
53,574-5	4-Amino-1-methyl-3-N-propyl-5-pyrazolecarboxamide, 96% [139756-02-8]	5g	97.20
★	(4-amino-2-methyl-5-propyl-2H-pyrazole-3-carboxylic acid amide) FW 182.22 mp 98-101° IRITANT		
27,018-0	3-(Aminomethyl)-PROXYL, free radical [54606-49-4] (3-aminomethyl-2,2,5,5-tetramethyl-1-pyrrolidinyloxy) FW 171.26 bp 128-131°/13mm Fp 144°F(62°C)	10mg	37.80
	Safety 2,194D R&S 1(2),2669B IRITANT	25mg	66.60
	Useful spin label for studying biological systems. Int. J. Pharm. 1993, 98, 131. Anal. Biochem. 1993, 212, 85.		
	4-Amino-N ¹⁰ -methylpteroylglutamic acid, see Amethopterin		
34,020-0	3-Amino-5-methylpyrazole, 97% [31230-17-8] FW 97.12 mp 45-47° bp 213°/14mm Fp >230°F(110°C) Beil. 25(4),2033 FT-NMR 1(3),70A R&S 1(2),2347A CORROSIVE	25g	51.00
		100g	141.20
A6,520-4	2-(Aminomethyl)pyridine, 99% [3731-51-9] (2-picolylamine) FW 108.14	5g	12.80
★	bp 82-85°/12mm n _D 1.5440 d 1.049 Fp 194°F(90°C) FT-NMR 1(3),288B FT-IR 1(2),764B Safety 2,195A R&S 1(2),2521E RTECS# US1840000 CORROSIVE	25g	42.20
		100g	118.40
A6,540-9	3-(Aminomethyl)pyridine, 99+% [3731-52-0] (3-picolylamine) FW 108.14 mp -21° bp 73-74°/11mm n _D 1.5510 d 1.062 Fp 213°F(100°C) FT-NMR 1(3),294A	5g	12.60
★	FT-IR 1(2),767D Safety 2,195B R&S 1(2),2525H IRITANT	100g	32.70
		500g	114.60
A6,560-3	4-(Aminomethyl)pyridine, 98% [3731-53-1] (4-picolylamine) FW 108.14 mp -8° bp 230° n _D 1.5510 d 1.065 Fp 227°F(108°C) Beil. 22(4),4181 FT-NMR 1(3),292A	5g	22.00
★	FT-IR 1(2),766C Safety 2,195C R&S 1(2),2523K IRITANT	25g	27.10
	Aminomethylpyridine, see Aminopicoline		
A6,570-0	2-Amino-4-methylpyrimidine, 97% [108-52-1] FW 109.13 mp 158-160° Beil. 24,84 FT-NMR 1(3),380C FT-IR 1(2),822C R&S 1(2),2575C RTECS# UV6485000 IRITANT	10g	34.10
		100g	250.30
	2-Amino-6-methyl-4-pyrimidinol, see A5,800-3, 2-Amino-4-hydroxy-6-methylpyrimidine. page 88		
42,288-6	(S)-(+)-2-(Aminomethyl)pyrrolidine, 97% [69500-64-7] FW 100.17 bp 65°/11mm n _D 1.4820 d 0.933 Fp 117°F(47°C) [α] _D +20° (c=1, CHCl ₃) Beil. 22(4),3761	100mg	26.70
	CORROSIVE MOISTURE-SENSITIVE	500mg	87.60
	Used in the synthesis of compounds with antitumor activity. J. Pharm. Sci. 1991, 80, 837.		
55,579-7	3-Amino-2-methyl-4(3H)-quinazolinone, 98% [1898-06-2] FW 175.19 mp 149-152°	5g	24.30
★	IRITANT	25g	78.80

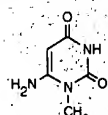


7-Amino-4-methyl-2(1H)-quinolinone, see 36,330-8, Carbostyrl 124 page 390		
(S)-(+)-2-Amino-4-(methylseleno)butanoic acid, see 47,394-4, L-(+)-Seleno methionine page 1643		
41,947-8	2-Amino-6-(methylsulfonyl)benzothiazole, 97% [17557-67-4] FW 228.29 mp 223°(dec.) Beil. 27(4),5473 TOXIC IRRITANT	100g 26.30
3-Aminomethyl-2,2,5,5-tetramethyl-1-pyrrolidinyloxy see 27,018-0, 3-(Amino-methyl)-PROXYL, free radical page 97		
A6,600-6	2-Amino-4-methylthiazole, 98% [1603-91-4] FW 114.17 mp 44-47° bp 231-232° Fp >230°F(110°C) Beil. 27,159 Merck Index 13,449 FT-NMR 1(3),114B FT-IR 1(2),645A R&S 1(2),2387G IRRITANT	5g 18.50 25g 61.30 100g 170.90
38,056-3	2-Amino-5-methylthiazole, 98% [7305-71-7] FW 114.17 mp 94-96° Beil. 27,162 FT-NMR 1(3),114C R&S 1(2),2387H IRRITANT	5g 16.70 25g 50.50
2-Amino-4-methylthio-1-butanol, see Methioninol		
3-Amino-4-methylthiophene-2-carboxylate, see 54,665-8, Methyl 3-amino-4-methylthiophene-2-carboxylate page 1210		
(1S,2S)-(+)-2-Amino-1-[4-(methylthio)phenyl]-1,3-propanediol, see 36,188-7		
(1S,2S)-(+)-Thiomcamine page 1775		
49,421-6	2-Amino-5-(methylthio)-1,3,4-thiadiazole, 99% [5319-77-7] FW 147.22 mp 178-181° Beil. 27(1),609	5g 27.00
19,068-3	3-Amino-5-methylthio-1H-1,2,4-triazole, 98% [45534-08-5] FW 130.17 mp 130-133° FT-NMR 1(3),90C FT-IR 1(2),629B Safety 2,196B R&S 1(2),2369A	25g 57.00
34,679-9	6-Amino-1-methyluracil, 97% [2434-53-9] FW 141.13 Beil. 24,470 FT-NMR 1(3),384C R&S 1(2),2579M	25g 45.50
A6,630-8	4-Aminomorpholine, 99% [4319-49-7] FW 102.14 bp 168° n _D 1.4770 d 1.059 Fp 137°F(58°C) Beil. 27,8 FT-NMR 1(1),600B FT-IR 1(1),384B Safety 2,196C R&S 1(1),403I RTECS# QD7300000 IRRITANT	10g 30.90 25g 61.50
A6,639-1	1-Aminonaphthalene [134-32-7], (α-naphthylamine, 1-naphthylamine) C ₁₀ H ₇ NH ₂ FW 143.19 mp 48-50° bp 301° d 1.114 Fp >230°F(110°C) Beil. 12,1212 Merck Index 13,6424 Safety 2,196D R&S 1(1),1445C RTECS# QM1400000 HIGHLY TOXIC CANCER SUSPECT AGENT OSHA-regulated carcinogen - 29 CFR Part 1910.1004	25g 36.20 250g 68.50
A6,640-5	2-Aminonaphthalene, 95% [91-59-8] (β-naphthylamine, 2-naphthylamine) C ₁₀ H ₇ NH ₂ FW 143.19 mp 111-113° bp 306° d 1.061 Beil. 12,1265 Merck Index 13,6425 Safety 2,197A R&S 1(1),1445J RTECS# QM2100000 HIGHLY TOXIC CANCER SUSPECT AGENT OSHA-regulated carcinogen - 29 CFR Part 1910.1009	1g 24.20
12,243-2	4-Amino-1-naphthalenecarbonitrile, 97% [58728-64-6] (4-cyano-1-naphthyl-amine) H ₂ NC ₁₀ H ₆ CN FW 168.20 mp 174.5-176.5° Beil. 14,533 FT-NMR 1(2),1549A FT-IR 1(2),464A Safety 2,197B R&S 1(2),2145L RTECS# QJ1865000	1g 45.50 5g 150.80
14,644-7	7-Amino-1,3-naphthalenedisulfonic acid, monopotassium salt hydrate, tech. 85% [303137-06-6] H ₂ NC ₁₀ H ₆ (SO ₃ H)SO ₃ K·xH ₂ O FW 341.41 mp >300° Beil. 14,784 Merck Index 13,412 FT-IR 1(2),501B Safety 2,197D R&S 1(2),2215H CORROSIVE	5g 16.10 100g 44.90
53,134-0	1-Amino-8-naphthalenesulfonic acid [82-75-7] (peri-acid) H ₂ NC ₁₀ H ₆ SO ₃ H FW 223.25 mp >350° Merck Index 13,6431 CORROSIVE	100g 42.50
29,113-7	2-Amino-1-naphthalenesulfonic acid, 98% [81-16-3] H ₂ NC ₁₀ H ₆ SO ₃ H FW 223.25 Beil. 14,736 Merck Index 13,6432 Safety 2,198A R&S 1(2),2199I RTECS# QK1250000 IRRITANT	5g 21.00 250g 38.50 1kg 103.50
25,061-9	4-Amino-1-naphthalenesulfonic acid, 97% [84-86-6] H ₂ NC ₁₀ H ₆ SO ₃ H FW 223.25 Beil. 14,739 Merck Index 13,6429 FT-IR 1(2),507C Safety 2,198B R&S 1(2),2213H RTECS# QK1270000 CORROSIVE	25g 69.00 100g 191.50
10,256-3	4-Amino-1-naphthalenesulfonic acid, sodium salt hydrate [123333-48-2] H ₂ NC ₁₀ H ₆ SO ₃ Na·xH ₂ O FW 245.23 mp 280°(dec.) Beil. 14,739 Merck Index 13,6323 FT-IR 1(2),507D Safety 2,198C R&S 1(2),2213I IRRITANT LIGHT-SENSITIVE	5g 14.10 100g 29.00 500g 88.80
28,507-2	5-Amino-2-naphthalenesulfonic acid, 95+% [119-79-9] H ₂ NC ₁₀ H ₆ SO ₃ H FW 223.25 Beil. 14,758 Merck Index 13,2372 Safety 2,198D R&S 1(2),2199J RTECS# QK1285000 CORROSIVE	5g 15.30 100g 51.30

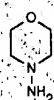


		US \$
4 page 390		
-Seleno		
28.29	100g	26.30
3-Amino-		
p 231-232°	5g	18.50
	25g	61.30
	100g	170.90
bell. 27,162	5g	16.70
	25g	50.50
3-amino-4-		
36,188-7		
22	5g	27.00
17	25g	57.00
2369A	25g	45.50
	10g	30.90
d 1.059	25g	61.50
196C	25g	36.20
ne) C ₁₀ H ₇ NH ₂	250g	68.50
Merck		
GHLY TOXIC		
lamine).....	1g	24.20
★		
GHLY TOXIC		
naphthyl-	1g	45.50
R 1(2),1549A	5g	150.80
hydrate, tech.	5g	16.10
10° Beil. 14,784	100g	44.90
CORROSIVE		
H ₂ SO ₄ H	100g	42.50
H ₂ H FW 223.25	5g	21.00
SH QK1250000	250g	38.50
	1kg	103.50
H ₂ H FW 223.25	25g	69.00
1(2),2213H	100g	191.50
333-48-2	5g	14.10
Index 13,6323	100g	29.00
ISITIVE	500g	88.80
SO ₃ H	5g	15.30
2199J	100g	51.30

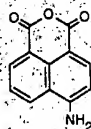
28,461-0	8-Amino-2-naphthalenesulfonic acid, 96% [119-28-8] H ₂ NC ₁₀ H ₆ SO ₃ H FW 223.25	5g	15.00
★	Beil. 14,765 Merck Index 13,2373 Safety 2,199A R&S 1(2),2199K CORROSIVE	25g	53.90
23,559-8	4-Amino-1,8-naphthalic anhydride, 95% [6492-86-0] FW 213.20 Beil. 18(2),469	250mg	31.50
★	FT-IR 1(2),336A Safety 2,199B R&S 1(2),1961B CANCER SUSPECT AGENT	1g	87.20
	MOISTURE-SENSITIVE		
	Contains 1-2% DMF		
A6,680-4	3-Amino-2-naphthoic acid, tech., 80% [5959-52-4] H ₂ NC ₁₀ H ₆ CO ₂ H FW 187.20	1g	20.50
★	mp 210°(dec.) Beil. 14,535 Merck Index 13,451 FT-NMR 1(2),1172B FT-IR 1(2),244A	5g	67.40
	R&S 1(2),1831N RTECS# QL1400000 IRRITANT		
51,237-0	6-Amino-2-naphthoic acid, 90% [116668-47-4] H ₂ NC ₁₀ H ₆ CO ₂ H FW 187.20	1g	33.90
★	mp 206-209° Beil. 14(2),324 IRRITANT	5g	111.60
	Contains ~5% 6-hydroxy-2-naphthoic acid and ~3% 5-amino-2-naphthoic acid		
13,347-7	1-Amino-2-naphthol hydrochloride, tech., 90% [1198-27-2] H ₂ NC ₁₀ H ₆ OH·HCl	1g	20.30
★	FW 195.65 mp 250°(dec.) Beil. 13,666 FT-IR 1(1),1252B R&S 1(1),1447B	5g	61.30
	RTECS# QL3335000 IRRITANT		
16,426-7	3-Amino-2-naphthol, 97% [5417-63-0] H ₂ NC ₁₀ H ₆ OH FW 159.19 mp 229-230°	1g	35.20
★	Beil. 13,681 FT-NMR 1(2),554A FT-IR 1(1),1253A Safety 2,199C R&S 1(1),1447G	5g	126.10
	IRRITANT		
13,348-5	4-Amino-1-naphthol hydrochloride, tech., 90% [5959-56-8] H ₂ NC ₁₀ H ₆ OH·HCl	5g	45.60
★	FW 195.65 mp 273°(dec.) Beil. 13,667 Merck Index 13,465 FT-IR 1(1),1252C	25g	176.00
	R&S 1(1),1447C RTECS# QL3342000 IRRITANT		
37,646-9	5-Amino-1-naphthol, 97% [83-55-6] H ₂ NC ₁₀ H ₆ OH FW 159.19 mp 190°(dec.)	25g	14.70
★	Beil. 13,670 FT-NMR 1(2),553C R&S 1(1),1447D IRRITANT	100g	33.60
53,691-1	8-Amino-2-naphthol, 97% [118-46-7] H ₂ NC ₁₀ H ₆ OH FW 159.19 mp 206-210°	25g	63.70
★	RTECS# QL3331000 IRRITANT LIGHT-SENSITIVE	500g	358.50
	Aminonaphtholsulfonic acid, see Aminohydroxynaphthalenesulfonic acid		
A6,820-3	6-Aminonicotinamide, 99% [329-89-5] FW 137.14 mp 245-248° FT-NMR 1(3),344A	250mg	25.60
★	FT-IR 1(2),799C Safety 2,200B R&S 1(2),2553A RTECS# US4550000 TERATOGEN	1g	66.30
		5g	220.00
A6,830-0	2-Aminonicotinic acid, 98% [5345-47-1] FW 138.13 mp 295-297°(dec.) Beil. 22,542	1g	11.10
★	FT-NMR 1(3),325B FT-IR 1(2),789D R&S 1(2),2543B IRRITANT	5g	28.90
		25g	122.70
21,687-9	6-Aminonicotinic acid, 97% [3167-49-5] FW 138.13 mp >300° Beil. 22,542	1g	29.80
★	Merck Index 13,455 FT-NMR 1(3),325C FT-IR 1(2),790A R&S 1(2),2543C	5g	96.40
41,893-5	2-(2-Amino-4-nitroanilino)ethanol, 99% [56932-44-6]	25g	20.20
★	H ₂ NC ₆ H ₃ (NO ₂)NHCH ₂ CH ₂ OH FW 197.19 mp 133-135° IRRITANT	100g	55.60
52,245-7	3-Amino-5-nitrobenzisothiazole, 97% [84987-89-3] FW 195.20 mp 250°(dec.)	5g	26.70
★	IRRITANT	100g	97.50
44,403-0	4'-Amino-5'-nitrobenzo-15-crown-5, 97% [77001-50-4] FW 328.32 mp 148-151°	250mg	73.00
★	IRRITANT	1g	208.60
	2-Aminonitrobenzoic acid, see Nitroanthranilic acid		
56,461-3	5-Amino-2-nitrobenzoic acid [13280-60-9] H ₂ NC ₆ H ₃ (NO ₂)CO ₂ H FW 182.14	1g	16.10
★	mp 236-238°(dec.)	5g	64.70
	2-Amino-5-nitrobenzonitrile, see 15,349-4, 5-Nitroanthranilonitrile page 1341		
32,469-8	4-Amino-3-nitrobenzonitrile, 98% [6393-40-4] H ₂ NC ₆ H ₃ (NO ₂)CN FW 163.14	1g	19.10
★	mp 161-163° Beil. 14(1),583 FT-NMR 1(2),1535C R&S 1(2),2137L IRRITANT	5g	68.70
21,173-7	2-Amino-5-nitrobenzophenone, 98+% [1775-95-7] H ₂ NC ₆ H ₃ (NO ₂)COC ₆ H ₅	25g	24.40
★	FW 242.23 mp 166-168° Beil. 14,79 FT-IR 1(2),73B Safety 2,200C R&S 1(2),1677G	100g	64.10
	RTECS# PC4935000 IRRITANT		
21,175-3	4-Amino-3-nitrobenzophenone, 98% [31431-19-3] H ₂ NC ₆ H ₃ (NO ₂)COC ₆ H ₅	5g	8.80
★	FW 242.23 mp 140-143° Beil. 14,86 FT-IR 1(2),73D Safety 2,200D R&S 1(2),1677I	25g	32.10
	IRRITANT		



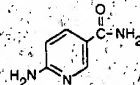
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A6,630-8



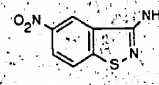
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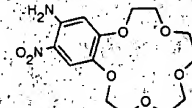
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A6,830-0



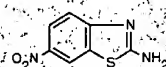
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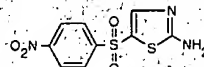
44,403-0

■ Aminonitro ■

			US \$
19,024-1	2-Amino-6-nitrobenzothiazole, 97+% [6285-57-0] FW 195.20 mp 247-249°	25g	20.60
★	Beil. 27(2); 232 FT-NMR 1(3); 209C FT-IR 1(2); 704D Safety 2,201A R&S 1(2); 2457J RTECS# DL0865000 IRRITANT	100g	48.90
	Aminonitrobenzotrifluoride, see Nitro(trifluoromethyl)aniline		
	2-Amino-6-[(4-nitrobenzyl)thio]-9-β-ribofuranosylpurine, see 86,166-9		
	S-(4-Nitrobenzyl)-6-thioguanosine page 1346		
A6,918-8	4-Amino-4'-nitrodiphenyl sulfide, 98% [101-59-7] [4-(4-nitrophenylthio)aniline]	5g	39.30
	H ₂ NC ₆ H ₄ SC ₆ H ₄ NO ₂ FW 246.29 mp 143-145° Beil. 13,534 FT-NMR 1(2); 780A FT-IR 1(1); 1386B R&S 1(1); 1603N RTECS# WP9640000 IRRITANT	25g	147.50
A7,000-3	1-Amino-4-nitronaphthalene, 97% [776-34-1] [4-nitro-1-naphthylamine]	1g	44.00
	H ₂ NC ₁₀ H ₆ NO ₂ FW 188.19 mp 190-193° Beil. 12,1259 FT-NMR 1(2); 784A FT-IR 1(1); 1390B Safety 2,202B R&S 1(1); 1607B RTECS# OM4370000 IRRITANT	5g	158.10
29,700-3	2-Amino-3-nitrophenol, 98% [603-85-0] [2-hydroxy-6-nitroaniline]	5g	50.70
	H ₂ NC ₆ H ₃ (NO ₂)OH FW 154.13 mp 212-213° Beil. 13(2); 191 FT-NMR 1(2); 713B Safety 2,202C R&S 1(1); 1567M IRRITANT	25g	164.50
A7,040-2	2-Amino-4-nitrophenol, 96% [99-57-0] [2-hydroxy-5-nitroaniline]	5g	25.50
★	H ₂ NC ₆ H ₃ (NO ₂)OH FW 154.13 mp 143-145° Beil. 13(2); 192 FT-NMR 1(2); 747A FT-IR 1(1); 1363A Safety 2,202D R&S 1(1); 1585H RTECS# SJ6300000 IRRITANT	25g	84.10
		100g	305.50
30,358-5	2-Amino-5-nitrophenol [121-88-0] [2-hydroxy-4-nitroaniline] H ₂ NC ₆ H ₃ (NO ₂)OH	25g	33.70
★	FW 154.13 mp 200°(dec.) Beil. 13,390 FT-NMR 1(2); 747B Safety 2,203A R&S 1(1); 1585I RTECS# SJ6302500 CANCER SUSPECT AGENT MUTAGEN	100g	93.60
A7,060-7	2-Amino-5-nitrophenol, tech.: 90% [121-88-0] [2-hydroxy-4-nitroaniline]	25g	13.80
★	H ₂ NC ₆ H ₃ (NO ₂)OH		
14,721-4	4-Amino-2-nitrophenol [119-34-6] [4-hydroxy-3-nitroaniline] H ₂ NC ₆ H ₃ (NO ₂)OH	25g	38.60
★	FW 154.13 mp 125-127° Beil. 13,520 FT-NMR 1(2); 747C FT-IR 1(1); 1363C Safety 2,203B R&S 1(1); 1585J RTECS# SJ6303000 CANCER SUSPECT AGENT MUTAGEN	100g	99.80
24,931-9	4-Amino-3-nitrophenol, 98% [610-81-1] H ₂ NC ₆ H ₃ (NO ₂)OH FW 154.13	5g	15.90
	mp 151-153° Beil. 13,521 FT-NMR 1(2); 748A R&S 1(1); 1585K IRRITANT	25g	59.30
		100g	164.40
A7,070-4	(1R,2R)-(-)-2-Amino-1-(4-nitrophenyl)-1,3-propanediol, 99% [716-61-0]	10g	53.20
	[p-(-)-threo] HOCH ₂ CH(NH ₂)CH(C ₆ H ₄ NO ₂)OH FW 212.21 mp 163-165° [α] _D ²⁵ -30° (c=1, 6N HCl) Beil. 13(3); 2267 FT-NMR 1(2); 705B FT-IR 1(1); 1344A Safety 2,203C R&S 1(1); 1563K RTECS# TY3100000 CORROSIVE	25g	95.00
47,167-4	(1S,2S)-(+)-2-Amino-1-(4-nitrophenyl)-1,3-propanediol, 99% [2964-48-9]	10g	20.50
	[L-(+)-threo-2-Amino-1-(4-nitrophenyl)-1,3-propanediol] O ₂ NC ₆ H ₄ CH(OH)CH(NH ₂)CH ₂ OH FW 212.21 mp 163-166° [α] _D ²³ +31° (c=1, 6N HCl) Beil. 13(3); 2267 CORROSIVE	50g	67.90
A7,075-5	2-Amino-5-(4-nitrophenylsulfonamido)thiazole [39565-05-4] FW 285.30 mp 222-226°	25g	20.50
	FT-IR 1(2); 648C R&S 1(2); 2391C RTECS# XJ2870000 IRRITANT	100g	53.30
11,351-4	2-Amino-3-nitropyridine, 99% [4214-75-9] FW 139.11 mp 165-167° Beil. 22(1); 631	5g	36.30
	FT-NMR 1(3); 305B FT-IR 1(2); 776A Safety 2,203D R&S 1(2); 2531L IRRITANT	25g	127.60
A7,080-1	2-Amino-5-nitropyridine, 97% [4214-76-0] FW 139.11 mp 186-188° Beil. 22(1); 631	5g	19.10
★	FT-NMR 1(3); 305C FT-IR 1(2); 776B Safety 2,204A R&S 1(2); 2531M IRRITANT		
A7,083-6	2-Amino-5-nitropyrimidine, 98% [3073-77-6] FW 140.10 mp 235-237°	250mg	24.10
	Beil. 24(1); 231 FT-IR 1(2); 833C R&S 1(2); 2583N IRRITANT	1g	68.60
		5g	255.90
14,502-5	5-Amino-6-nitroquinoline, 97% [35975-00-9] FW 189.17 mp 272-273°(dec.)	1g	18.30
	FT-IR 1(2); 868C R&S 1(2); 2627G IRRITANT	5g	66.90
13,350-7	2-Amino-5-nitrothiazole, 97% [121-66-4] FW 145.14 mp 202°(dec.)	25g	22.10
★	Merck Index 13,456 FT-IR 1(2); 647C Safety 2,204B R&S 1(2); 2389C RTECS# XJ2800000 CANCER SUSPECT AGENT	100g	55.40
	2-Amino-6-nitrotoluene, see 11,584-3, 2-Methyl-3-nitroaniline page 1261		
	3-Aminonoradamantane hydrochloride, see 29,187-0, 3-Noradamantanamine hydrochloride page 1371		



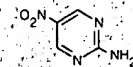
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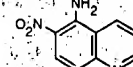
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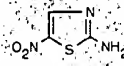
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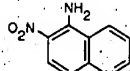
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47-249°	25g	20.60
1(2),2457J	100g	48.90
5,166-9		
ylthio)aniline]	5g	39.30
),780A	25g	147.50
imine)	1g	44.00
4A	5g	158.10
IRITANT		
)	5g	50.70
2),713B	25g	164.50
	5g	25.50
2),747A	25g	84.10
IRITANT	100g	305.50
gH ₃ (NO ₂)OH ...	25g	33.70
IA	100g	93.60
ITAGEN		
iniline)	25g	13.80
gH ₃ (NO ₂)OH	25g	38.60
63C	100g	99.80
AGENT		
13	5g	15.90
IT	25g	59.30
	100g	164.40
-61-0]	10g	53.20
5°	25g	95.00
1344A		
64-48-9]	10g	20.50
(c=1, 6N HCl)	50g	67.90
0 mp 222-226°	25g	20.50
	100g	53.30
Beil. 22(1),631	5g	36.30
ITANT	25g	127.60
Beil. 22(1),631	5g	19.10
ITANT		
37°	250mg	24.10
	1g	68.60
	5g	255.90
73°(dec.)	1g	18.30
	5g	66.90
	25g	22.10
	100g	55.40

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nantanamine



17,083-6

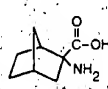


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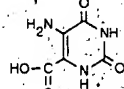
13,351-5	2-Aminonorbomane hydrochloride, 99% [14370-45-7] (2-norbomanamine hydrochloride) FW 147.65 mp 295°(dec.) [α] _D ²⁵ 0° (c=1, H ₂ O) Beil. 12(3),160. FT-IR 1(1),320C R&S 1(1),343H. Predominantly endo	1g	35.90
17,960-4	exo-2-Aminonorbomane, 99% [7242-92-4] (exo-2-norbomanamine) FW 111.19 bp 49°/10mm n _D ²⁰ 1.4800 d 0.938 Fp 95°F(35°C) [α] _D ²⁵ 0° (neat) Beil. 12(3),160. FT-NMR 1(1),516C FT-IR 1(1),320B Safety 2,204C R&S 1(1),343G. FLAMMABLE LIQUID IRRITANT	1g	37.80
22,552-5	2-Amino-2-norbomanecarboxylic acid, 98%, mixture of isomers [20448-79-7] (BCH) FW 155.20 mp >300° FT-NMR 1(1),885B FT-IR 1(1),585A R&S 1(1),663G. Inhibitor of amino acid transport across cell membranes. J. Biol. Chem. 1992, 267, 25951. Am. J. Physiol. 1992, 262, R761; Chem. Abstr. 1992, 117, 64436p. Aminooctanoic acid, see Aminocaprylic acid	100mg	38.80
		250mg	61.20
		1g	167.60
19,121-3	5-Aminoorotic acid, 99% [7164-43-4] (5-amino-2,6-dioxo-1,2,3,6-tetrahydro-4-pyrimidinecarboxylic acid) FW 171.11 mp >300° Beil. 25,264 FT-NMR 1(3),390C FT-IR 1(2),835B R&S 1(2),2585J RTECS# RM3190000 IRRITANT	50g	33.60
37,179-3	2-Amino-4-oxo-4H-1-benzopyran-3-carboxaldehyde, see 40,202-8, 2-Amino-3-formylchromone page 86		
	5-(3-Amino-5-oxo-2-pyrazolin-1-yl)-2-phenoxybenzenesulfonic acid, tech., 92% [30479-81-3] FW 347.35 mp 297°(dec.) R&S 1(2),2199B IRRITANT	5g	18.40
	(Aminooxy)acetic acid hemihydrochloride, see C1,340-8, Carboxymethoxylamine hemihydrochloride page 392	25g	60.60
A7,090-9	(+)-6-Aminopenicillanic acid, 96% [551-16-6] (6-APA) FW 216.26 mp 198-200°(dec.) [α] _D ²² +276.3° (c=1.2, 0.1N HCl) Merck Index 13,457 FT-IR 1(1),791C Safety 2,204D R&S 1(1),927C RTECS# XH8225000	10g	29.70
		50g	106.40
	3-Aminopentane, see 19,019-5, 1-Ethylpropylamine page 882		
	2-Aminopentanoic acid, see Norvaline		
23,668-3	D,L-2-Amino-1-pentanol, 97% [4146-04-7] CH ₃ CH ₂ CH ₂ CH(NH ₂)CH ₂ OH FW 103.17 bp 194-195° n _D ²⁰ 1.4510 d 0.922 Fp 203°F(95°C) Beil. 4(3),794 FT-NMR 1(1),540A FT-IR 1(1),338B Safety 2,205A R&S 1(1),363I. IRRITANT. HYGROSCOPIC	1g	34.40
		10g	203.60
53,457-9	(R)-(-)-2-Amino-1-pentanol [80696-30-6] (D-norvalinol) CH ₃ (CH ₂) ₂ CH(NH ₂)CH ₂ OH FW 103.17 mp 44-48° Fp 204°F(95°C) [α] _D ²⁵ -17° (c=1, CHCl ₃) IRRITANT	1g	69.40
53,458-7	(S)-(+)-2-Amino-1-pentanol [22724-81-8] (L-norvalinol) CH ₃ (CH ₂) ₂ CH(NH ₂)CH ₂ OH FW 103.17 mp 44-48° Fp 214°F(101°C) [α] _D ²⁵ +17° (c=1, CHCl ₃) IRRITANT	1g	69.40
12,304-8	5-Amino-1-pentanol, 95% [2508-29-4] H ₂ N(CH ₂) ₄ OH FW 103.17 mp 35-37° bp 122°/16mm n _D ²⁰ 1.4610 d 0.949 Fp 150°F(65°C) Beil. 4(1),441 FT-NMR 1(1),539C FT-IR 1(1),339A Safety 2,205B R&S 1(1),363H IRRITANT	10g	18.40
		50g	60.90
17,344-4	D,L-2-Amino-4-pentenoic acid, 99% [7685-44-1] H ₂ C=CHCH ₂ CH(NH ₂)CO ₂ H FW 115.13 mp 258-260° Beil. 4,467 FT-NMR 1(1),874C FT-IR 1(1),574D Safety 2,205C R&S 1(1),655N IRRITANT	1g	19.30
		5g	74.20
28,501-3	L-2-Amino-4-pentenoic acid, 98% [16338-48-0] H ₂ C=CHCH ₂ CH(NH ₂)CO ₂ H FW 115.13 mp 283°(dec.) [α] _D ²⁵ -36° (c=4, H ₂ O) Beil. 4(4),2852 FT-NMR 1(1),875A Safety 2,205D R&S 1(1),655O IRRITANT	100mg	31.70
		1g	182.10
23,212-2	2-Aminoperimidine hydrobromide hydrate, 98% [313223-13-1] FW 264.13 mp 297° FT-IR 1(2),893B Safety 2,206A LIGHT-SENSITIVE	5g	101.30
		25g	337.40
14,910-1	9-Aminophenanthrene, 96% [947-73-9] (9-phenanthrenamine) FW 193.25 mp 137-139° Beil. 12,1338 FT-IR 1(1),1255C R&S 1(1),1451B RTECS# SG0175000 IRRITANT	250mg	136.80
		1g	356.20
	4-Aminophenazone, see A3,930-0, 4-Aminoantipyrine page 68		



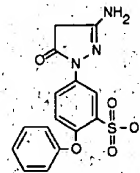
13,351-5



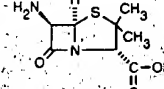
22,552-5



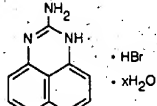
19,121-3



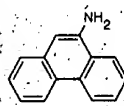
37,179-3



A7,090-9



23,212-2

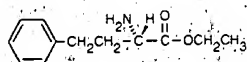


14,910-1

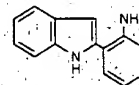
OH FW 137.18.....	5g	104.60
(3), 1679	25g	320.80
OH FW 137.18.....	5g	26.40
Safety 2.206B	25g	84.60
ylamine page 103		
4-177° Beil. 13.354 ..	5g	17.60
2.206C	100g	19.40
	500g	62.50
24-126°	5g	16.90
B FT-IR 1(1), 1200A	100g	18.10
ITANT	500g	49.50
FW 116.19.....	1g	283.00
188-190°	5g	16.70
10B Safety 2.207A	100g	16.90
	500g	51.90
	1kg	78.30
	250g	28.10
	1kg	83.90
HCl FW 145.59.....	5g	15.00
1(1), 1399G	100g	22.70
	500g	94.60
O ₂ H FW 151.17.....	1g	47.30
IRITANT	10g	313.10
H FW 151.17.....	5g	14.90
2A' FT-IR 1(2), 163B	25g	56.40
	100g	144.70
anide page 71		
anide page 71		
	250mg	29.00
+37° (c=2, H ₂ O)	1g	79.90
	1g	36.10
) Beil. 14(4), 1688	5g	127.60
	250mg	19.50
) Beil. 14(4), 1687	1g	47.20
[312693-79-1].....	250mg	28.60
It hydrate, tech.	1g	28.40
max 404nm	10g	136.70
2-0].....	5g	83.30
x 461nm IRRITANT		
1-1', Mordant Yellow		
1-1] (3-amino-.....	1g	51.00
-96° Beil. 16(3), 1284	5g	163.40
inobenzene.....	5g	118.20
00- FT-IR 1(2), 1136C	25g	390.70
-1].....	1g	46.40
2- HCl FW 173.41.....	5g	153.70

ax 1-800-962-9591

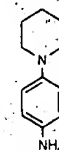
2-Amino-4-phenylbutyric acid, see Homophenylalanine		
33,533-9 4-(4-Aminophenyl)butyric acid, 95% [15118-60-2] H ₂ NC ₆ H ₄ (CH ₂) ₃ CO ₂ H.....	1g	20.40
FW 179.22 mp 121-124° Beil. 14(3), 1279 FT-NMR 1(2), 1022C R&S 1(2), 1759M	5g	55.60
IRRITANT		
53,291-6 (S)-(+)-2-Amino-4-phenylbutyric acid ethyl ester hydrochloride, 97%.....	1g	53.90
C ₁₆ H ₁₉ (CH ₂) ₂ CH(NH ₂)CO ₂ C ₂ H ₅ ·HCl FW 243.73 mp 159-163° [α] _D ²⁵ +26° (c=1, CHCl ₃)	5g	141.70
16,676-6 2-Aminophenyl disulfide, 97% [1141-88-4] (2,2'-dithiodianiline) (H ₂ NC ₆ H ₄) ₂ S ₂	10g	51.40
FW 248.37 mp 91-92° Beil. 13.400 FT-NMR 1(2), 465B FT-IR 1(1), 1197A Safety 2.208B	50g	156.60
R&S 1(1), 1389D RTECS# BX9540000 IRRITANT		
36,946-2 4-Aminophenyl disulfide, 98% [722-27-0] (H ₂ NC ₆ H ₄) ₂ S ₂ FW 248.37 mp 77-78°.....	5g	39.10
Beil. 13.536 FT-NMR 1(2), 486B R&S 1(1), 1399J RTECS# BX9580000 IRRITANT	25g	129.40
1-(3-Aminophenyl)ethanol, see H ₂ 620-0, 3-(1-Hydroxyethyl)aniline page 1026		
54,438-8 1-(4-Aminophenyl)ethanol, 97% [14572-89-5] H ₂ NC ₆ H ₄ CH(OH)CH ₃ FW 137.18.....	5g	52.10
mp 70-74° IRRITANT	25g	173.50
49,457-7 (R)-(-)-2-Amino-1-phenylethanol, 90% [2549-14-6] H ₂ NCH ₂ CH(C ₆ H ₅)OH.....	1g	85.30
FW 137.18 mp 59-65° Fp >230°F (110°C) [α] _D ²⁵ -39° (c=2, C ₂ H ₅ OH) Beil. 13(4), 1801		
CORROSIVE		
May contain up to 10% water		
A7,240-5 2-Amino-1-phenylethanol, 98% [7568-93-6] [(α-aminomethyl)benzylalcohol].....	10g	55.80
H ₂ NCH ₂ CH(C ₆ H ₅)OH FW 137.18 mp 56-58° bp 160°/17mm Merck Index 13.7371	50g	185.40
FT-NMR 1(2), 575C FT-IR 1(1), 1272B Safety 2.208D R&S 1(1), 1469G		
RTECS# DN5500000 IRRITANT		
2-Amino-2-phenylethanol, see 2-Phenylglycinol		
Aminophenyl ether, see Oxydianiline		
12,305-6 2-(4-Aminophenyl)ethylamine, 97% [13472-00-9] (4-aminophenethylamine).....	1g	31.20
H ₂ NC ₆ H ₄ CH ₂ CH ₂ NH ₂ FW 136.20 bp 103°/0.3mm n _D ²⁰ 1.5910 d 1.034	5g	103.40
Fp >230°F (110°C) Beil. 13(1), 48 FT-NMR 1(2), 611C FT-IR 1(1), 1289B Safety 2.209B	10g	198.10
R&S 1(1), 1489H CORROSIVE		
Reagent used in polycondensation reactions. <i>Macromolecules</i> 1994, 27, 1289. <i>Polym. J.</i> (Tokyo) 1991, 23, 1511; <i>Chem. Abstr.</i> 1992, 116, 106932j		
3-(4-Aminophenyl)-3-ethyl-2,6-piperidinedione, see 25,919-5. Aminogluthethimide		
page 86		
1-(4-Aminophenyl)-1H-imidazole, see 44,439-1, 4-(1H-imidazol-1-yl)aniline		
page 1060		
37,686-8 2-(2-Aminophenyl)indole, 97% [32566-01-1] FW 208.26 mp 154-155°.....	250mg	44.60
Beil. 22(4), 5017 FT-NMR 1(3), 133A R&S 1(2), 2405O IRRITANT	1g	116.80
10,556-2 4-Aminophenylmercuric acetate, tech. [6283-24-5] [4-(acetoxymethyl)aniline].....	5g	20.60
CH ₃ CO ₂ HgC ₆ H ₄ NH ₂ FW 351.76 mp 163-165° FT-IR 1(2), 274D Safety 2.209D	25g	78.50
R&S 1(2), 1867B RTECS# OV5550000 HIGHLY TOXIC IRRITANT		
54,827-8 N-(3-Aminophenyl)methanesulfonamide [37045-73-1] FW 186.23 mp 117-121°.....	5g	26.10
IRRITANT	25g	86.80
37,177-7 N-(4-Aminophenyl)-4-methylbenzenesulfonamide, 99% [6380-08-1].....	1g	15.30
CH ₃ C ₆ H ₄ SO ₂ NHC ₆ H ₄ NH ₂ FW 262.33 mp 183-185° Beil. 13.114 FT-NMR 1(2), 1635C	10g	78.20
R&S 1(2), 2241I IRRITANT		
39,411-4 2-Amino-4-phenylphenol, tech., 90% [1134-36-7] H ₂ NC ₆ H ₃ (C ₆ H ₅)OH FW 185.23.....	1g	21.20
mp 198-202° Beil. 13(2), 419 IRRITANT	5g	63.90
Aminophenyl phenyl ether, see Phenoxylaniline		
2-Aminophenyl phenyl sulfone, see 22,504-5, 2-(Phenylsulfonyl)aniline		
page 1462		
55,662-9 N-(4-Aminophenyl)piperidine, 97% [2359-60-6] FW 176.26 mp 26-29° n _D ²⁰ 1.5937.....	1g	20.20
Fp >230°F (110°C) IRRITANT	5g	67.20
55,673-4 N-(3-Aminophenyl)propanamide, 97% [22987-10-6] H ₂ NC ₆ H ₄ NHCOC ₂ H ₅	25g	69.40
FW 164.20 mp 92-96° IRRITANT		
1-Amino-2-phenylpropane, see 18,007-6, β-Methylphenethylamine page 1272		



53,291-6



37,686-8

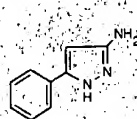


55,662-9

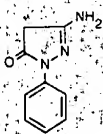
C of A's / MSDS's and spectra now available online at www.sigma-aldrich.com.

Aminopheny

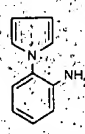
		US \$
24,888-6	(1R,2R)-(-)-2-Amino-1-phenyl-1,3-propanediol, 98% [46032-98-8] C ₆ H ₅ CH(OH)CH(NH ₂)CH ₂ OH FW 167.21 mp 112-118° [α] _D ²⁵ -37° (c=1, 1N HCl) Beil 13(4), 2697 Fieser 13.18 R&S 1(1), 1475J CORROSIVE	1g 36.00 5g 107.50
18,654-6	(1S,2S)-(+)-2-Amino-1-phenyl-1,3-propanediol, 98% [28143-91-1] C ₆ H ₅ CH(OH)CH(NH ₂)CH ₂ OH FW 167.21 mp 109-113° [α] _D ²⁵ +37° (c=1, 1N HCl) Beil 13(4), 2698 Fieser 12.32 13.18 FT-NMR 1(2), 587C FT-IR 1(1), 1278B Safety 2, 211A R&S 1(1), 1475I CORROSIVE Precursor to chiral 2-oxazolines. J. Am. Chem. Soc. 1976, 98, 567	1g 23.20 5g 71.60
28,449-1	(R)-(+)-2-Amino-3-phenyl-1-propanol, 98% [5267-64-1] (D-phenylalaninol) C ₆ H ₅ CH ₂ CH(NH ₂)CH ₂ OH FW 151.21 mp 93-95° [α] _D ²⁵ +23° (c=1.2, 1N HCl) FT-NMR 1(2), 571B R&S 1(1), 1465L CORROSIVE	1g 31.70 5g 117.10
09222	(S)-2-Amino-1-phenylethanol, purum, ≥97.0% (GC, sum of enantiomers) [56613-81-1] [(S)-α-(aminomethyl)benzylalcohol] FW 137.18	100mg 58.70 500mg 232.10
19,043-8	(S)-(-)-2-Amino-3-phenyl-1-propanol, 98% [3182-95-4] (L-phenylalaninol) C ₆ H ₅ CH ₂ CH(NH ₂)CH ₂ OH FW 151.21 mp 92-94° [α] _D ²² -22.8° (c=1.2, 1N HCl) Beil 13(3), 1757 FT-NMR 1(2), 576C FT-IR 1(1), 1272D Safety 2, 211B R&S 1(1), 1469J RTECS# UA6900000 IRRITANT Reacts with nitriles to form oxazolines ¹⁻³ which are useful in Pd-catalyzed allylic substitution ^{2,3} Also employed in amidation for chiral resolution ⁴ and NADH modeling. (1) Tetrahedron Lett. 1993, 34, 7725. (2) ibid. 1993, 34, 3149. (3) ibid. 1993, 34, 2015. (4) ibid. 1993, 34, 7081. (5) Tetrahedron 1993, 49, 5237.	1g 23.10 10g 152.60
15,949-2	D,L-3-Amino-3-phenylpropionic acid, 98% [614-19-7] (β-aminohydrocinnamic acid) C ₆ H ₅ CH(NH ₂)CH ₂ CO ₂ H FW 165.19 mp 222°(dec.) Beil. 14,493 FT-NMR 1(2), 995C FT-IR 1(2), 146B R&S 1(2), 1747D	1g 14.10 5g 46.70 25g 178.50
56,025-1	3-(4-Aminophenyl)propionic acid, 97% [2393-17-1] H ₂ NC ₆ H ₄ (CH ₂) ₂ CO ₂ H FW 165.19 mp 133-137° IRRITANT	1g 34.40 5g 142.50
39,379-7	3-Amino-5-phenylpyrazole, 98% [1572-10-7] FW 159.19 mp 124-127° Beil. 24,148 RTECS# UQ6109000 IRRITANT	1g 18.80 5g 62.20
A7,400-9	3-Amino-1-phenyl-2-pyrazolin-5-one, 99+% [4149-06-8] FW 175.19 mp 210-215°(dec.) FT-NMR 1(3), 71B FT-IR 1(2), 608C R&S 1(2), 2347E RTECS# URO393500 IRRITANT	25g 18.90 100g 53.10
19,694-0	1-(2-Aminophenyl)pyrrole, 98+% [6025-60-1] FW 158.20 mp 96-98° FT-NMR 1(3), 4B FT-IR 1(2), 567A Safety 2, 211C R&S 1(2), 2303K IRRITANT	1g 20.30 10g 112.60
A7,460-2	3-Aminophenyl sulfone, 97% [599-61-1] (3,3'-sulfonyldianiline) (H ₂ NC ₆ H ₄) ₂ SO ₂ FW 248.30 mp 170-173° Beil. 13,426 FT-IR 1(2), 490C Safety 2, 211D R&S 1(2), 2191B RTECS# BY8800000 IRRITANT	25g 17.30 100g 48.60
A7,480-7	4-Aminophenyl sulfone, 97% [80-08-0] (dapsone, DDS, 4,4'-sulfonyldianiline) (H ₂ NC ₆ H ₄) ₂ SO ₂ FW 248.30 mp 175-177° Beil. 13,536 Merck Index 13,2847 FT-IR 1(2), 490B Safety 2, 212A R&S 1(2), 2191A RTECS# BY8925000 IRRITANT	5g 13.30 100g 30.70 500g 109.70
36,152-6	2-(3-Aminophenylsulfonyl)ethanol hydrochloride, 97% [19076-03-0] H ₂ NC ₆ H ₄ SO ₂ CH ₂ CH ₂ OH, HCl FW 237.71 mp 210°(dec.) Beil. 13,426 FT-NMR 1(2), 1584A R&S 1(2), 2187C IRRITANT	1g 70.30 5g 234.20
14,105-4	2-Amino-4-phenyl-5-tetradecylthiazole, 98% [64415-14-1] FW 372.62 mp 72-74° FT-IR 1(2), 646B R&S 1(2), 2387L	5g 47.80
46,462-7	2-Amino-5-phenyl-1,3,4-thiadiazole sulfate, 97% [312619-47-9] FW 177.23 mp 200°(dec.) IRRITANT	1g 13.50 5g 44.20
A7,500-5	2-Amino-4-phenylthiazole hydrobromide monohydrate, 99% [52253-69-7] FW 275.17 mp 180-183° Beil. 27,204 FT-NMR 1(3), 115B FT-IR 1(2), 646A R&S 1(2), 2387K	25g 73.80 100g 176.20
43,855-3	4-Amino-5-phenyl-4H-1,2,4-triazole-3-thiol, 97% [21089-45-2] FW 192.24 mp 195°(dec.) Beil. 26(4), 472 IRRITANT	1g 13.50 5g 44.20



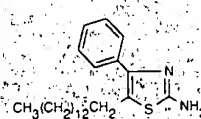
39,379-7



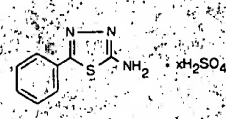
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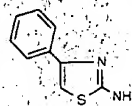
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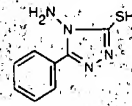
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46,462-7

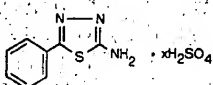


A7,500-5



43,855-3

	1g	US \$ 36.00
(1N HCl)	5g	107.50
	1g	23.20
(1N HCl)	5g	71.60
788		
alaninol)	1g	31.70
HCl)	5g	117.10
omers)	100mg	58.70
	500mg	232.10
alaninol)	1g	23.10
NHCl)	10g	152.60
S 1(1), 1469J		
d allylic		
H modeling ⁵		
93, 34, 2015		
drocinnamic	1g	14.10
	5g	46.70
	25g	178.50
2)CO ₂ H	1g	34.40
	5g	142.50
27° Beil. 24,148	1g	18.80
	5g	62.20
9	25g	18.90
	100g	53.10
	1g	20.30
ANT	10g	112.60
H ₂ NC ₆ H ₄ SO ₂	25g	17.30
S 1(2), 2191B	100g	48.60
onyldianiline)	5g	13.30
2847	100g	30.70
IRITANT	500g	109.70
3-0)	1g	70.30
	5g	234.20
62 mp 72-74°	5g	47.80
W 177.23	1g	13.50
	5g	44.20
253-69-7)	25g	73.80
3A	100g	176.20
192.24	1g	13.50
	5g	44.20

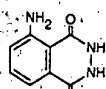


46,462-7

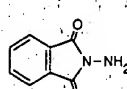


855-3

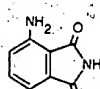
12,307-2	3-Aminophthalhydrazide, 97% [521-31-3] (5-amino-2,3-dihydro-1,4-phthal- azinedione, luminol) FW 177.16 mp >300° Beil. 25(1),698 Merck Index 13,5622 FT-IR 1(2),384D Safety 2,213C R&S 1(2),2041F RTECS# TH8890060 IRRITANT Used for chemiluminescence analysis ^{1,2} of, for example, metal cations ³ and blood. ⁴ (1) Endeavour 1992, 16, 194. (2) Aldrichimica Acta 1983, 16, 59. (3) Bull. Chem. Soc. Jpn. 1992, 65, 1591. (4) J. Biochem. Biophys. Methods 1993, 27, 301.	2.5g	US \$ 17.20
39,123-9	3-Aminophthalhydrazide, sodium salt hemihydrate, 98% [206658-90-4] FW 199.15 mp >300° Beil. 25(1),698 IRRITANT	1g	29.90
19,317-8	4-Aminophthalhydrazide hydrate, 98% [3682-14-2] (6-amino-2,3-dihydro-1,4- phthalazinedione) FW 177.16 mp 300° Beil. 25,487 FT-IR 1(2),385A R&S 1(2),2041G IRRITANT Produces chemiluminescence. Aldrichimica Acta 1983, 16, 59.	5g	116.70
33,534-7	3-Aminophthallic acid, tech., 90% [5434-20-8] H ₂ NC ₆ H ₃ 1,2-(CO ₂ H) ₂ FW 181.15 mp 180-185° Beil. 14,552 R&S 1(2),1799K IRRITANT	250mg	18.90
52,471-6	4-Aminophthallic acid, 97% [5434-21-9] H ₂ NC ₆ H ₃ 1,2-(CO ₂ H) ₂ FW 181.15 mp 344°(dec.) IRRITANT	1g	41.20
17,831-4	N-Aminophthalimide, tech., 90% [1875-48-5] FW 162.15 mp 200-202° Fieser 1,38 5,14 FT-IR 1(2),407B R&S 1(2),2083F IRRITANT Employed in the azidination of chiral N-enoyl sulfamates. J. Chem. Soc., Chem. Commun. 1993, 1074. Remainder phthalhydrazide	1g	22.90
52,479-4	3-Aminophthalimide, 97% [2518-24-3] FW 162.15 mp 268.5-273.5° IRRITANT	5g	78.30
33,601-7	4-Aminophthalonitrile, 98% [56765-79-8] H ₂ NC ₆ H ₃ 1,2-(CN) ₂ FW 143.15 mp 179-181° Beil. 14(3),1399	10g	130.00
A7,563-3	2-Amino-3-picoline, 95% [1603-40-3] (2-amino-3-methylpyridine) FW 108.14 mp 32-34° bp 221-222° n _D 1.5820 d 1.073 Fp 233°F(111°C) Beil. 22(2),342 FT-NMR 1(3),296A FT-IR 1(2),769A Safety 2,214A R&S 1(2),2527C RTECS# US1850000 HIGHLY TOXIC IRRITANT	1g	23.60
12,308-0	2-Amino-4-picoline, 99% [695-34-1] (2-amino-4-methylpyridine) FW 108.14 mp 98-100° bp 230° Beil. 22(2),342 Merck Index 13,465 FT-NMR 1(3),298C FT-IR 1(2),771B Safety 2,214B R&S 1(2),2527L RTECS# TJ5150000 HIGHLY TOXIC IRRITANT	5g	78.20
A7,568-4	2-Amino-5-picoline, 99% [1603-41-4] (2-amino-5-methylpyridine, 6-amino-3- picoline) FW 108.14 mp 76-77° bp 227° FT-NMR 1(3),298B FT-IR 1(2),771A Safety 2,214C R&S 1(2),2527K RTECS# TJ5141000 HIGHLY TOXIC IRRITANT	10g	130.20
A7,570-6	2-Amino-6-picoline, 98% [1824-81-3] (2-amino-6-methylpyridine, 6-amino-2- picoline) FW 108.14 mp 41.5-45.5° bp 208-209° Fp 218°F(103°C) Beil. 22(1),633 FT-NMR 1(3),295C FT-IR 1(2),768D Safety 2,214D R&S 1(2),2527B RTECS# US1885000 HIGHLY TOXIC IRRITANT	5g	39.80
A7,590-0	4-Amino-1-piperazineethanol, see 12,296-3; 1-Amino-4-(2-hydroxyethyl)- piperazine page 88	25g	132.10
29,336-9	3-Aminopiperidine dihydrochloride, 97% [138060-07-8] FW 173.09 mp 230°(dec.) [α] _D ²⁰ 0° (c=1, H ₂ O) Beil. 22(4),3743 Safety 2,215B R&S 1(1),389D IRRITANT	5g	14.90
53,666-0	(S)-(+)-3-Aminopiperidine dihydrochloride, 97% FW 173.08 mp 212°(dec.) [α] _D ²⁵ +2° (c=1, H ₂ O) IRRITANT	10g	59.00
56,147-9	4-Aminopiperidine [13035-19-3] FW 100.17 n _D 1.4910 d 0.954 Fp 114°F(45°C) CORROSIVE HYGROSCOPIC	25g	122.00
56,385-4	4-Amino-2-piperidin-1-yl-pyrimidine-5-carbonitrile, 97% [90973-23-2] FW 203.24 mp 224° IRRITANT	100g	355.00
	2-Amino-1,3-propanediol, see Serinol	50mg	36.70
		250mg	126.30
		1g	96.40
		5g	60.60
		25g	202.00
		1g	55.60



12,307-2



17,831-4



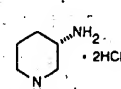
52,479-4



A7,563-3



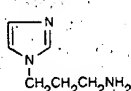
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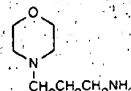
53,666-0

H ₂ O FW 91.11	5g	11.20
0° (c=7.3, H ₂ O)	25g	37.90
TY2800000	100g	107.60
CH(OH)CH ₂ OH	1g	30.90
°C [η] ²³ -12° (neat)		
2) ₃ SO ₃ H FW 139.17	1g	25.80
	5g	69.70
lar to those elicited by		
em. Abstr. 1977, 86,		
59, 373.		
V(CH ₂) ₂ SO ₃ Na · xH ₂ O	1g	46.60
R&S 1(1), 1059H	5g	155.00
mine)	1g	35.80
0.954 Fp 165°F(73°C)	5g	141.50
7.1(1), 336A		
1) ₂ CH(OH)CH ₂ NH ₂	5mL	19.20
18° (neat) Beil. 4,289	100mL	21.30
IM	1L	25.90
	2L	46.40
imine)	1g	34.40
0.954 Fp 169°F(76°C)	5g	136.10
1(1), 335D		
CH ₂ CH(NH ₂)CH ₂ OH	1g	24.70
-18° (neat)	5g	74.50
CH(NH ₂)CH ₂ OH	5g	38.60
4(1), 432 Merck	25g	134.30
R&S 1(1), 361J		
CH ₂ CH(NH ₂)CH ₂ OH	1g	25.00
α) ²⁵ +18° (neat)	10g	168.60
1,335B Safety 2,217A		
d-catalyzed allylic		
1, 34, 3149, (3)		
11 mp 10-12°	10g	16.00
NMR 1(1), 536A	50g	56.30
000 CORROSIVE		
	100g	15.20
	500g	34.60
	6x500g	134.30
chlorotriyl 3-amino-		
HO(CH ₂) ₃ NH ₂	10mL	24.90
SIVE		
-amino-1,1,3-tri-	5g	23.10
FW 132.13		
1), 861C Safety 2,217C		
	5g	27.10
FT-NMR 1(1), 1381A	25g	115.10
00000 TERATOGEN		

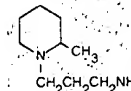
- 36,425-8 3-Aminopropyl-functionalized silica gel, suitable for low- and medium-pressure reversed-phase liquid chromatography R&S 1(3), 3569A IRRITANT
Useful as a heterogeneous, reusable catalyst for the Knoevenagel condensation. J. Chem. Soc., Perkin Trans. 1 1989, 105.
-1 meq NH₂/g
-9% functionalized
- 27,226-4 1-(3-Aminopropyl)imidazole, 98% [5036-48-6] FW 125.18 n_D²⁰ 1.5190 d 1.049
Fp >230°F(110°C) Beil. 23(4), 577 FT-NMR 1(3), 81B R&S 1(2), 2353N CORROSIVE
- 42,211-8 O-(2-Aminopropyl)-O-(2-methoxyethyl)polypropylene glycol [77110-54-4]
CH₃OCH₂CH₂(OC₃H₆)_nOCH₂CH₂(NH₂)CH₃ n_D²⁰ 1.4450 d 0.980 Fp >230°F(110°C)
TOXIC IRRITANT
Average M_n ca. 600
- 12,309-9 4-(3-Aminopropyl)morpholine, 98% [123-00-2] FW 144.22 mp -15° bp 224°
n_D²⁰ 1.4760 d 0.987 Fp 210°F(98°C) FT-NMR 1(1), 601A FT-IR 1(1), 384D Safety 2,218B
R&S 1(1), 403L RTECS# QD7700000 CORROSIVE
- 47,188-7 (1*R*)-(-)-(1-Aminopropyl)phosphonic acid, 98% [98049-00-4]
C₂H₅CH(NH₂)P(O)(OH)₂ FW 139.09 mp 265-269° [α]_D²⁵ -16.5° (c=2, 1N NaOH)
IRRITANT
- 47,189-5 (1*S*)-(+)-(1-Aminopropyl)phosphonic acid, 98% [98048-99-8]
C₂H₅CH(NH₂)P(O)(OH)₂ FW 139.09 mp 265-269° [α]_D²⁵ +16.5° (c=2, 1N NaOH)
IRRITANT
- 26,861-5 3-Aminopropylphosphonic acid, 99% [13138-33-5] H₂N(CH₂)₃P(O)(OH)₂
FW 139.09 mp 294°(dec.) Beil. 4(3), 1793 FT-NMR 1(1), 1477C Safety 2,218D
R&S 1(1), 1103B IRRITANT
- 18,611-2 1-(3-Aminopropyl)-2-pipecoline, 96% [25560-00-3] FW 156.27 bp 96-97°/15mm
n_D²⁰ 1.4760 d 0.889 Fp 191°F(88°C) FT-NMR 1(1), 576A FT-IR 1(1), 370A Safety 2,219A
R&S 1(1), 389C CORROSIVE
- 1-100-6 N-(3-Aminopropyl)-1,3-propanediamine, 98% [56-18-8] HN(CH₂)₃NH₂
FW 131.22 mp -14° bp 151°/50mm n_D²⁰ 1.4810 d 0.938 Fp 245°F(118°C)
FT-NMR 1(1), 500C FT-IR 1(1), 310B Safety 2,1975D R&S 1(1), 331E
RTECS# JL9450000 HIGHLY TOXIC CORROSIVE
- 13,656-5 1-(3-Aminopropyl)-2-pyrrolidinone, tech. [7663-77-6] FW 142.20
bp 120-123°/1mm n_D²⁰ 1.5000 d 1.014 Fp >230°F(110°C) FT-NMR 1(1), 1289B
FT-IR 1(1), 790D Safety 2,219B R&S 1(1), 925D RTECS# UY5739500 CORROSIVE
Contains ~2% 2-pyrrolidinone and varying amounts of 1,5-diazabicyclo[4.3.0]non-5-ene
(Aldrich cat. no. 13,658-1)
- 44,014-0 3-Aminopropyltriethoxysilane, 99% [919-30-2] [3-(triethoxysilyl)propylamine]
H₂N(CH₂)₃Si(OC₂H₅)₃ FW 221.37 bp 217° d 0.946 Fp 205°F(96°C)
RTECS# TX2100000 CORROSIVE MOISTURE-SENSITIVE
DOW CORNING[®] product Z-6011
(Packaged under nitrogen in Sure/Seal[™] bottles)
- 28,177-8 3-Aminopropyltrimethoxysilane, 97% [13822-56-5] [3-(trimethoxysilyl)propylamine]
H₂N(CH₂)₃Si(OCH₃)₃ FW 179.29 bp 91-92°/15mm n_D²⁰ 1.4240 d 1.027
Fp 197°F(91°C) Safety 2,219D R&S 1(2), 2979J, CORROSIVE MOISTURE-SENSITIVE
- 2-Amino-4-pteridinol, see 28,708-3, Pterin page 1601
- 86,227-4 Aminopterine hydrate, 97% [54-62-6] (4-aminofolic acid, 4-aminopteroyl-
glutamic acid) FW 440.42 mp 225°(dec.) [α]_D²⁴ +18° (c=0.1, 0.1N NaOH)
Beil. 26(4), 831 Merck Index 13,471 FT-NMR 1(3), 481A R&S 1(2), 2663E
RTECS# MA1050000 HIGHLY TOXIC TERATOGEN
- 4-Aminopteroylglutamic acid, see 86,227-4, Aminopterine hydrate page 107
- 28,508-0 2-Aminopurine [452-06-2] FW 135.13 mp 280-282° Beil. 26,414 R&S 1(2), 2471I
RTECS# UO7475000
- 6-Aminopurine, see Adenine



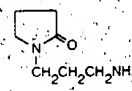
27,226-4



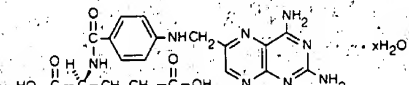
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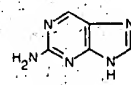
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13,656-5



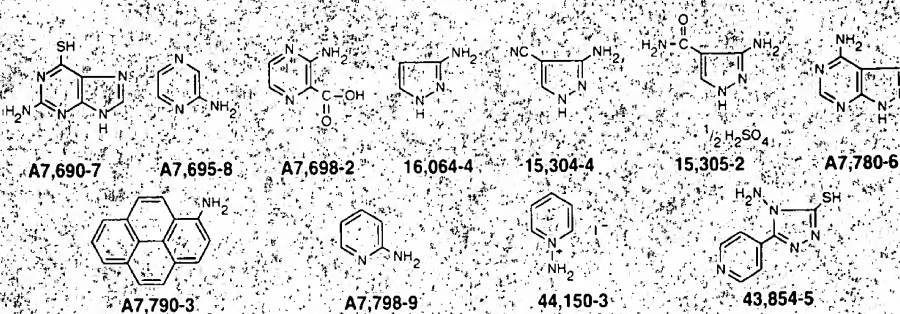
44,014-0



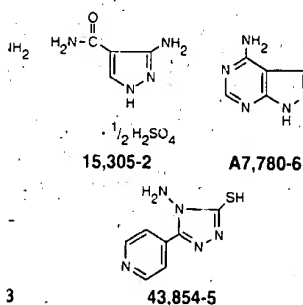
28,177-8

■ Aminopurin ■

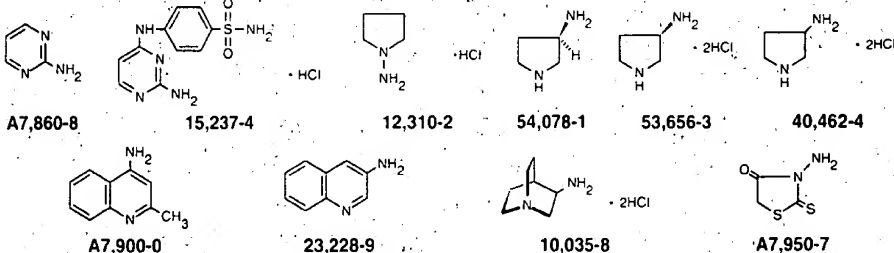
		US \$
A7,690-7	2-Amino-6-purinethiol, 98% [154-42-7] (6-thioguanine) FW 167.19 mp >360° Merck Index 13,9411 FT-IR 1(2),715D Safety 2,220B R&S 1(2),2471K RTECS# UP0740000 TOXIC	1g 47.80 5g 200.80
A7,695-8	Aminopyrazine, 99% [5049-61-6] FW 95.11 mp 119-120° Beil: 24,82 FT-NMR 1(3),406C FT-IR 1(2),843C R&S 1(2),2593J IRRITANT	1g 11.40 10g 59.60 50g 214.80
A7,698-2	3-Aminopyrazine-2-carboxylic acid, 99+% [5424-01-1] FW 139.11 mp 205°(dec.) FT-NMR 1(3),407B FT-IR 1(2),843D Safety 2,220C R&S 1(2),2593M IRRITANT	1g 16.40 5g 54.30 25g 185.80
16,064-4	3-Aminopyrazole, 98% [1820-80-0] FW 83.09 mp 37-39° bp 218°/122mm Fp >230°F(110°C) Beil: 24,14 FT-NMR 1(3),69C FT-IR 1(2),607D Safety 2,220D R&S 1(2),2345N CORROSIVE Used in the preparation of heterocyclic compounds of pharmaceutical interest. Trends, Heterocycl. Chem. 1991, 2, 97	2.5g 45.10 10g 156.10
15,304-4	3-Amino-4-pyrazolecarbonitrile, 97% [16617-46-2] (3-amino-4-cyanopyrazole) FW 108.10 mp 172-174° FT-NMR 1(3),75A FT-IR 1(2),611A Safety 2,221A R&S 1(2),2349E CORROSIVE	1g 31.60 5g 107.00 25g 221.70
15,305-2	3-Amino-4-pyrazolecarboxamide hemisulfate [27511-79-1] FW 175.16 mp 224°(dec.) FT-IR 1(2),610D R&S 1(2),2349D	1g 27.00 5g 105.30
A7,740-7	3-Amino-4-pyrazolecarboxylic acid, 95% [41680-34-6] FW 127.10 mp 135°(dec.) FT-NMR 1(3),73A FT-IR 1(2),609C Safety 2,221B R&S 1(2),2347K RTECS# UQ6390000 IRRITANT	1g 74.50 5g 203.10
A7,780-6	4-Aminopyrazolo[3,4-d]pyrimidine, 98% [20289-44-5] FW 135.13 mp >325° FT-IR 1(2),707A R&S 1(2),2463G RTECS# UR0717000 TOXIC IRRITANT Decreases serum cholesterol markedly in rats. Science 1976, 193, 903. Biochemistry 1984, 23, 4533	250mg 16.00 1g 41.20 5g 190.40
	4-Aminopyrazolo[3,4-d]pyrimidine-6-thiol, see A6,020-2; 4-Amino-6- mercaptopyrazolo[3,4-d]pyrimidine, page 91	
A7,790-3	1-Aminopyrene, 97% [1606-67-3] (1-pyrenamine) FW 217.27 mp 115-117° Beil: 12,1341 FT-IR 1(1),1256B Safety 2,221C R&S 1(1),1451E RTECS# UR2275000	250mg 15.40 1g 42.30
A7,798-9	2-Aminopyridine, 99+% [504-29-0] FW 94.12 mp 59-60° bp 204-210° Fp 198°F(92°C) Beil: 22,428 Fieser 9,18 Merck Index 13,472 FT-IR 1(2),761B Safety 2,221D R&S 1(2),2519B RTECS# US1575000 HIGHLY TOXIC FLAMMABLE SOLID	25g 27.30 100g 64.70
A7,799-7	2-Aminopyridine, 99% [504-29-0]	5g 10.80 100g 23.50 500g 88.20
A7,820-9	3-Aminopyridine, 99% [462-08-8] FW 94.12 mp 57-60° bp 248° Beil: 22,431 Merck Index 13,472 FT-NMR 1(3),293C FT-IR 1(2),767C Safety 2,222A R&S 1(2),2525G RTECS# US1650000 HIGHLY TOXIC IRRITANT	25g 16.40 100g 38.50 500g 133.60
27,587-5	4-Aminopyridine, 99+% [504-24-5] FW 94.12 mp 160-162° bp 273° Beil: 22,433 Merck Index 13,3964 FT-NMR 1(3),291A FT-IR 1(1),765D Safety 2,222B R&S 1(2),2523H RTECS# US1750000 HIGHLY TOXIC IRRITANT	1g 14.40 5g 46.90
A7,840-3	4-Aminopyridine, 98% [504-24-5]	25g 20.90 100g 71.00
44,150-3	1-Aminopyridinium iodide, 97% [6295-87-0] FW 222.03 mp 159-161° IRRITANT	1g 9.80 10g 53.90
	2-Amino-3-pyridinol, see 12,251-3; 2-Amino-3-hydroxypyridine, page 89	
43,854-5	4-Amino-5-(4-pyridyl)-4H-1,2,4-triazole-3-thiol, 97% [36209-51-5] FW 193.23 mp 250-254° Beil: 26(4),2129 IRRITANT	1g 16.00 10g 87.90



67.19 mp -360°	1g	US \$ 47.80
71K	5g	200.80
24.R2	1g	11.40
	10g	59.60
	50g	214.80
39.11 mp 205°(dec.)	1g	16.40
M-IRRITANT	5g	54.30
	25g	185.80
8"/122mm	2.5g	45.10
Safety 2.220D	10g	156.10
Interest, Trends		
4-cyanopyrazole	1g	31.60
y 2.221A	5g	107.00
	25g	221.70
W 175.16	1g	27.00
	5g	105.30
27.10 mp 135°(dec.)	1g	74.50
	5g	203.10
5.13 mp >325°	250mg	16.00
IRRITANT	1g	41.20
903. Biochemistry	5g	190.40
mino-6-		
mp 115-117°	250mg	15.40
ECSS# UR2275000	1g	42.30
4-210°	25g	27.30
IR 1(2).761B	100g	64.70
KIC		
	5g	10.80
	100g	23.50
	500g	88.20
Beil. 22.431	25g	16.40
222A	100g	38.50
	500g	133.60
273° Beil. 22.433	1g	14.40
:222B	5g	46.90
	25g	20.90
	100g	71.00
59-161° IRRITANT...	1g	9.80
	10g	53.90
page 89		
79-51-5] FW 193.23	1g	16.00
	10g	87.90

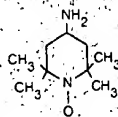


A7,860-8	2-Aminopyrimidine, 97% [109-12-6] FW 95.11 mp 125-127° Beil. 24.80	5g	US \$ 8.20
★	FT-NMR 1(3).380A FT-IR 1(2).822B Safety 2.222D R&S 1(2).2573M	100g	19.50
	RTECS# UVG326000 IRRITANT	500g	71.60
26,182-3	4-Aminopyrimidine, 98% [591-54-8] FW 95.11 mp 154-156° Beil. 24.81	250mg	41.20
	FT-NMR 1(3).380B Safety 2.223A R&S 1(2).2575B IRRITANT	1g	122.50
	Aminopyrimidinediol, see Aminodihydroxypyrimidine		
	4-Amino-N-(2-pyrimidinyl)benzenesulfonamide, see 28,719-9, Sulfadiazine page 1699		
15,237-4	N-(2-Amino-4-pyrimidinyl)sulfanilamide monohydrochloride, 98%	100g	16.60
✓	[22199-93-5] FW 301.76 mp 284-286°(dec.) FT-NMR 1(3).393B Safety 2.223B		
	R&S 1(2).2587H IRRITANT		
	Aminopyrine, see D13,910-6, 4-(Dimethylamino)antipyrene page 711		
12,310-2	1-Aminopyrrolidine hydrochloride, 97% [63234-71-9] FW 122.60 mp 117-119°	5g	140.90
✓	FT-IR 1(1).357D Safety 2.223C R&S 1(1).379J RTECS# UX9693000 HYGROSCOPIC	10g	233.20
54,078-1	(R)-(+)-3-Aminopyrrolidine, 98% [116183-82-5] FW 86.14 bp 164-165° n _D 1.4890	1g	79.30
✓	d 0.984 Fp 147°F(63°C) IRRITANT		
54,080-3	(S)-(-)-3-Aminopyrrolidine, 98% [128345-57-3] FW 86.14 bp 164-165° n _D 1.4880	1g	90.50
✓	d 0.967 Fp 155°F(68°C) [α] _D -20° (neat) IRRITANT HYGROSCOPIC		
53,656-3	(R)-(+)-3-Aminopyrrolidine dihydrochloride, 97% [116183-81-4] FW 159.06	1g	90.80
✓	mp 340°(dec.) [α] _D -1.8° (c=1, H ₂ O) IRRITANT HYGROSCOPIC		
40,462-4	3-Aminopyrrolidine dihydrochloride, 98% [103831-11-4] FW 159.06 mp >300°	5g	45.20
✓	[α] _D 0° (c=1, H ₂ O) IRRITANT HYGROSCOPIC	25g	154.90
A7,900-0	4-Aminoquinoline, 98% [6628-04-2] FW 158.20 mp 167-169° bp 333° Beil. 22.453	5g	12.70
✓	FT-NMR 1(3).437C FT-IR 1(2).863C R&S 1(2).2621D IRRITANT	25g	42.50
23,228-9	3-Aminoquinoline, 98% [580-17-6] (3-quinolinamine) FW 144.18 mp 91-92°	1g	12.70
✓	Beil. 22(1).638 FT-NMR 1(3).436B FT-IR 1(2).862D Safety 2.223D R&S 1(2).2619L	5g	43.60
★	RTECS# VA9622000 IRRITANT		
A7,920-5	5-Aminoquinoline, 97% [611-34-7] (5-quinolinamine) FW 144.18 mp 107-109°	1g	40.40
	bp 310° Beil. 22.445 FT-NMR 1(3).436C FT-IR 1(2).863B R&S 1(2).2621A	5g	143.80
	RTECS# VA9625000 IRRITANT		
27,558-1	6-Aminoquinoline, 98% [580-15-4] (6-quinolinamine) FW 144.18 mp 117-119°	1g	28.90
	bp 146°/0.3mm Beil. 22.447 FT-NMR 1(3).437A Safety 2.224A R&S 1(2).2621B	5g	93.80
	IRRITANT		
	Fluorescent derivatizing agent useful in the detection of biochemicals, Anal. Biochem.		
	1993, 211, 279, J. Chromatogr. 1992, 600, 279		
26,078-9	8-Aminoquinoline, 98% [578-66-5] (8-quinolinamine) FW 144.18 mp 66-68°	1g	11.00
★	bp 174°/26mm Beil. 22.450 FT-NMR 1(3).437B Safety 2.224B R&S 1(2).2621C	5g	33.70
	RTECS# VA9627000 IRRITANT	25g	133.10
	5-Amino-8-quinolinol dihydrochloride, see 30,552-9, 5-Amino-8-hydroxy-quinoline dihydrochloride page 89		
41,571-5	(R)-(+)-3-Aminoquinuclidine dihydrochloride, 98% [123536-14-1] FW 199.14	250mg	19.60
	mp >300° [α] _D +24° (c=1, H ₂ O) RTECS# VD4590000 IRRITANT	1g	54.20
10,035-8	3-Aminoquinuclidine dihydrochloride, 98% [6530-09-2] FW 199.14	1g	22.10
✓	mp 321-323°(dec.) [α] _D 24.0° (c=1, CH ₃ OH) Beil. 22(4).3816 FT-IR 1(1).380C	5g	73.20
	R&S 1(1).399E RTECS# VD4590000	25g	233.10
41,572-3	(S)-(-)-3-Aminoquinuclidine dihydrochloride, 98% [119904-90-4] FW 199.14	250mg	18.10
	[α] _D -24° (c=1, H ₂ O) RTECS# VD4590000 IRRITANT HYGROSCOPIC	1g	49.20
16,317-1	4-Aminoresorcinol hydrochloride, 97% [34781-86-7] H ₂ NC ₆ H ₃ .1.3-(OH) ₂ .HCl	1g	38.90
	FW 161.59 mp 220°(dec.) Beil. 13.783 FT-IR 1(1).1226A Safety 2.224C	5g	128.50
	R&S 1(1).1415G LIGHT-SENSITIVE		
A7,950-7	3-Aminorhodanine, 99% [1438-16-0] FW 148.21 mp 100-103° Beil. 27.245	10g	39.60
★	FT-IR 1(1).835C R&S 1(1).979L HIGHLY TOXIC		



■ Aminosalicylic ■

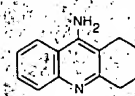
			US \$
25,530-0	3-Aminosalicylic acid, 97% [570-23-0] $H_2NC_6H_3-2-(OH)CO_2H$ FW 153.14 mp 240°(dec.) Beil. 14,577 FT-NMR 1(2), 1099A Safety 2,224D R&S 1(2), 1799D IRRITANT	100mg 500mg	44.00 184.60
A7,960-4	4-Aminosalicylic acid, 99% [65-49-6] $H_2NC_6H_3-2-(OH)CO_2H$ FW 153.14 mp 147°(dec.) Beil. 14,579 Merck Index 13,476 FT-NMR 1(2), 1122A FT-IR 1(2), 219C Safety 2,225A R&S 1(2), 1809C RTECS# VO1225000 IRRITANT LIGHT-SENSITIVE	5g 100g 500g	8.50 15.00 58.50
85,654-1	4-Aminosalicylic acid, sodium salt dihydrate, 99% [6018-19-5] $H_2NC_6H_3-2-(OH)CO_2Na \cdot 2H_2O$ FW 211.15 mp 250-254°(dec.) Merck Index 13,476 FT-IR 1(2), 275A R&S 1(2), 1867C IRRITANT	5g 100g 1kg	7.90 20.60 102.70
A7,980-9	5-Aminosalicylic acid, 95% [89-57-6] $H_2NC_6H_3-2-(OH)CO_2H$ FW 153.14 mp 280°(dec.) Beil. 14,579 Merck Index 13,5931 FT-NMR 1(2), 1121C FT-IR 1(2), 219B Safety 2,225B R&S 1(2), 1809A RTECS# VO1400000 IRRITANT	5g 100g 500g	12.20 33.50 107.30
	4-Aminostyrene, see 4-Vinylaniline		
	2-Aminosuccinamic acid, see Asparagine		
	Aminosuccinic acid, see Aspartic acid		
	Aminosulfo-2-Ar functionalized silica gel, see 53,793-4, 4-Ethyl-benzene-sulfonamide-Functionalized silica gel page 839		
	(S)-(-)-5-(Aminosulfonyl)-N-[(1-ethyl-2-pyrrolidinyl)methyl]-2-methoxybenzamide, see 44,559-2, (S)-(-)-Sulpiride, page 1707		
16,394-5	4-Amino-TEMPO, free radical, 97% [14691-88-4] (4-amino-2,2,6,6-tetramethylpiperidinyloxy, free radical) FW 171.26 FT-IR 1(2), 908C R&S 1(2), 2673L Useful spin label for studying biological systems ¹ and polymers. ² (1) Biochemistry 1994, 33, 3855; (2) J. Colloid Interface Sci., 1994, 165, 236	250mg 1g 5g	25.20 72.10 235.30
38,107-1	2-Aminoterephthalic acid, 99% [10312-55-7] $H_2NC_6H_3-1,4-(CO_2H)_2$ FW 181.15 mp 324°(dec.) Beil. 14,558 FT-NMR 1(2), 1134B R&S 1(2), 1815B IRRITANT	25g	26.40
10,061-7	4-Amino-2,3,5,6-tetrafluorobenzamide, 97% [1548-74-9] $H_2NC_6F_4CONH_2$ FW 208.11 mp 185-186° FT-NMR 1(2), 1402A FT-IR 1(2), 375B Safety 2,225C R&S 1(2), 2023L IRRITANT	5g	34.60
38,170-5	2-Amino-3,4,5,6-tetrafluorobenzoic acid, 99% [1765-42-0] (3,4,5,6-tetrafluoroanthranilic acid) $H_2NC_6F_4CO_2H$ FW 209.10 mp 141-143° IRRITANT	250mg 1g	27.70 75.80
24,793-6	4-Amino-2,3,5,6-tetrafluorobenzoic acid, 99% [944-43-4] $H_2NC_6F_4CO_2H$ FW 209.10 mp 185-187° FT-NMR 1(2), 1158C FT-IR 1(2), 235C Safety 2,225D R&S 1(2), 1825H IRRITANT	1g 5g	28.80 110.10
24,794-4	4-Amino-2,3,5,6-tetrafluorobenzonitrile, 99% [17823-38-0] $H_2NC_6F_4CN$ FW 190.10 mp 96-98° FT-IR 1(2), 461D Safety 2,226A R&S 1(2), 2143A IRRITANT	1g 5g	18.50 57.90
30,062-4	4-Amino-2,3,5,6-tetrafluoropyridine, 98% [1682-20-8] FW 166.08 mp 85-87° FT-NMR 1(3), 301C Safety 2,226B R&S 1(2), 2529K IRRITANT	1g	37.30
A7,992-2	9-Amino-1,2,3,4-tetrahydroacridine hydrochloride hydrate, 99+% (tetrahydroaminacrine hydrochloride, THA hydrochloride) FW 234.73 mp 284-286° Beil. 22(4), 4911 Merck Index 13,9003 FT-NMR 1(3), 443B Safety 2,226C R&S 1(2), 2625K RTECS# AR9532100 HIGHLY TOXIC IRRITANT	1g 5g	20.00 62.50
09440	(R)-(+)-3-Aminotetrahydrofuran toluene-4-sulfonate, puriss., ≥99.0% (T) [111769-27-8] [(R)-(+)-3-aminotetrahydrofuran p-toluenesulfonate] FW 259.33 mp 139-141° $[\alpha]_D^{25} +4.5 \pm 0.5^\circ$ (c=3 in MeOH) Chiral amine; amidation with amino acids. K. Barros et al., Liebig's Ann. Chem. 1986, 1950.	250mg 1g	45.80 158.50
	Aminotetrahydronaphthalene, see Tetrahydronaphthylamine		
	2-Amino-3,4,5,6-tetrahydropyridine hydrochloride, see 13,117-2, 2-Imino-piperidine hydrochloride, page 1061		
	1-Aminotetralin, see 1,2,3,4-Tetrahydro-1-naphthylamine		
11,573-8	4-Amino-2,2,6,6-tetramethylpiperidine, 98% [36768-62-4] FW 156.27 mp 16-18° bp 188-189° $n_D^{20} 1.4700$ d 0.912 Fp 162°F (72°C) Beil. 22(1), 627 FT-NMR 1(1), 577C FT-IR 1(1), 369C Safety 2,226D R&S 1(1), 389I RTECS# TM4290100 IRRITANT	5g 25g	29.80 128.60
	4-Amino-2,2,6,6-tetramethylpiperidinyloxy, free radical, see 16,394-5		
55,072-8	5-Aminotetrazole, 97% [4418-61-5] FW 85.07 mp 201-205° RTECS# XF7465000	100g 500g	23.00 76.40



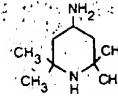
16,394-5



30,062-4

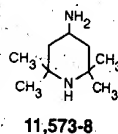


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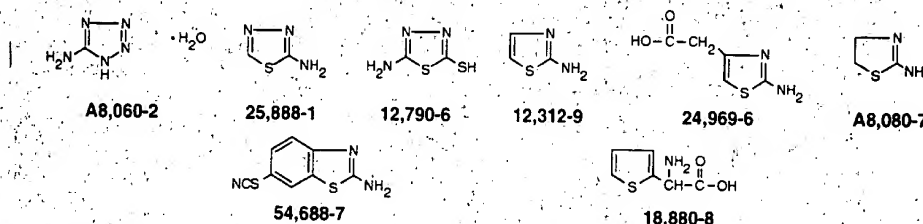


11,573-8

W 153.14	100mg	44.00
1(2),1799D	500mg	184.60
V 153.14	5g	8.50
FT-IR 1(2),219C	100g	15.00
GHT-SENSITIVE	500g	58.50
5)	5g	7.90
* Index 13,476	100g	20.60
	1kg	102.70
V 153.14	5g	12.20
C FT-IR 1(2),219B	100g	33.50
	500g	107.30
hyl-benzene-		
nethoxy		
6,6-tetramethyl-	250mg	25.20
),2673L	1g	72.10
Biochemistry 1994,	5g	235.30
2H ₂ FW 181.15	25g	26.40
RITANT		
6F ₄ CONH ₂	5g	34.60
ty 2,225C		
1,5,6-tetrafluoro-	250mg	27.70
NT	1g	75.80
C ₆ F ₄ CO ₂ H	1g	28.80
ty 2,225D	5g	110.10
VC ₆ F ₄ CN	1g	18.50
I3A IRRITANT	5g	57.90
08 mp 85-87°	1g	37.30
9+%	1g	20.00
/ 234.73	5g	62.50
3B Safety 2,226C		
99.0% (T)	250mg	45.80
nate]FW 259.33	1g	158.50
1. Chem. 1986,		
7-2, 2-Imino-		
156.27 mp 16-18°	5g	29.80
NMR 1(1),577C	25g	128.60
IRRITANT		
6,394-5,		
CS# XF7465000	100g	23.00
	500g	76.40



A8,060-2	5-Aminotetrazole monohydrate, 97% [15454-54-3] FW 103.08 mp 200-204°(dec.)	5g	12.50
★	Beil 26,403 FT-IR 1(2),633D Safety 2,227A R&S 1(2),2373D RTECS# XF7465000	100g	32.50
	FLAMMABLE SOLID	500g	135.00
	Useful reagent in condensation reactions to form, for example, pyrimidines. J. Heterocycl. Chem. 1993, 30, 1267. Indian J. Chem., Sect. B 1993, 32B, 886.		
25,888-1	2-Amino-1,3,4-thiadiazole, 97% [4005-51-0] FW 101.13 mp 190-192° Beil. 27,624	1g	10.70
★	FT-IR 1(2),650C Safety 2,227B R&S 1(2),2393E RTECS# XI3000000 TOXIC	5g	23.80
	IRRITANT		
12,790-6	5-Amino-1,3,4-thiadiazole-2-thiol, 98% [2349-67-9] FW 133.20 mp 235°(dec.)	10g	15.90
★	Beil 27,674 FT-IR 1(2),651C R&S 1(2),2393I RTECS# XI4550000 IRRITANT	50g	45.70
12,312-9	2-Aminothiazole, 97% [96-50-4] FW 100.14 mp 91-93° Beil. 27,155 Fieser 4,19	5g	15.40
★	Merck Index 13,479 FT-NMR 1(3),114A FT-IR 1(2),644D Safety 2,227C	100g	21.10
	R&S 1(2),2387F RTECS# XJ2100000 TOXIC	500g	69.30
24,969-6	2-Amino-4-thiazoleacetic acid, 95% [29676-71-9] FW 158.18 mp 130°(dec.)	25g	38.70
★	Beil 27,336 FT-IR 1(2),647D Safety 2,227D R&S 1(2),2389F IRRITANT	100g	107.50
A8,080-7	2-Amino-2-thiazoline, 97% [1779-81-3] FW 102.16 mp 79-82° Beil. 27,136	5g	11.30
★	FT-NMR 1(1),1333B FT-IR 1(1),824B R&S 1(1),969B RTECS# XJ6540000	25g	25.90
		100g	64.60
26,372-9	2-Amino-2-thiazoline hydrochloride, 98% [3882-98-2] FW 138.62 mp 200-202°	250g	38.10
★	Beil 27,136 R&S 1(1),969C HYGROSCOPIC	1kg	94.90
54,688-7	2-Amino-6-thiocyanatobenzothiazole, 96% [7170-77-6] FW 207.27 mp 194-197°	25g	164.90
★	IRRITANT		
18,880-8	α-Amino-2-thiopheneacetic acid, 97% [21124-40-3] FW 157.19 mp 208-210°	1g	49.90
★	Beil 18,631 FT-IR 1(2),603D R&S 1(2),2341E	5g	174.30
	α-Amino-2-thiophenepropionic acid, see 3-(2-Thienyl)alanine		
27,424-0	2-Aminothiophenol, 99% [137-07-5] (2-aminobenzenethiol) H ₂ NC ₆ H ₄ SH	25g	29.40
★	FW 125.19 mp 19-21° bp 70-72°/0.2mm n _D ²⁰ 1.6420 d 1.170 Fp 175°F(79°C)	100g	97.20
	Beil 13,397 FT-NMR 1(2),464C Safety 2,228B R&S 1(1),1389B RTECS# DC0600000		
	CORROSIVE STENCH		
12,313-7	2-Aminothiophenol, tech., 90% [137-07-5] (2-aminobenzenethiol) H ₂ NC ₆ H ₄ SH	100g	28.80
★		500g	77.80
14,348-0	3-Aminothiophenol, 96% [22948-02-3] (3-aminobenzenethiol) H ₂ NC ₆ H ₄ SH	5g	75.00
	FW 125.19 n _D ²⁰ 1.6560 d 1.179 Fp 175°F(79°C) Beil. 13,425 FT-NMR 1(2),471A	25g	251.60
	FT-IR 1(1),1200C Safety 2,228C R&S 1(1),1391I CORROSIVE STENCH		
42,296-7	4-Aminothiophenol, 97% [1193-02-8] NH ₂ C ₆ H ₄ SH FW 125.19 mp 39-42°	5g	32.10
	bp 140-145°/16mm Fp >230°F(110°C) RTECS# DC0602500 CORROSIVE STENCH	25g	106.90
A8,100-5	4-Aminothiophenol, tech., 90% [1193-02-8] (4-aminobenzenethiol) H ₂ NC ₆ H ₄ SH	1g	10.90
		5g	22.00
	6-Amino-2-thiouracil monohydrate, see A5,740-6, 4-Amino-6-hydroxy-2-mercaptopyrimidine monohydrate page 88		
	α-Amino-p-toluenesulfonamide hydrochloride, see A6,180-2, 4-(Aminomethyl)-benzenesulfonamide hydrochloride page 93		
10,255-5	2-Aminotoluene-5-sulfonic acid, 98% [98-33-9] (4-amino-3-methyl-benzenesulfonic acid) H ₂ NC ₆ H ₃ (CH ₃)SO ₃ H FW 187.22 mp >300° Beil. 14,726	250g	14.40
★	CANCER SUSPECT AGENT MUTAGEN		
	DuPont product		
56,292-0	2-Aminotoluene-5-sulfonic acid, 97% [98-33-9] (4-amino-3-methyl-benzenesulfonic acid) H ₂ NC ₆ H ₃ (CH ₃)SO ₃ H FW 187.22 mp 290°(dec.)	250g	13.90
★	CANCER SUSPECT AGENT CORROSIVE	1kg	38.20
21,876-6	4-Aminotoluene-3-sulfonic acid, 99% [88-44-8] (2-amino-5-methyl-benzenesulfonic acid) H ₂ NC ₆ H ₃ (CH ₃)SO ₃ H FW 187.22 Beil. 14,723	250g	28.20
★	FT-NMR 1(2),1595C FT-IR 1(2),495B Safety 2,188A R&S 1(2),2197E		
	RTECS# XT6320000 CANCER SUSPECT AGENT CORROSIVE		
	DuPont product		



■ Aminotolyl ■

			US \$
53,416-1	2-Amino-4-(p-tolyl)-thiazole, 96% [2103-91-5] (4-(4-methylphenyl)-2-thiazolamine) FW 190.26 mp 132-136°	1g	12.60
10,062-5	3-Amino-1,2,4-triazine, 97% [1120-99-6] FW 96.09 mp 174-177° FT-IR 1(2).846D R&S 1(2).2599E RTECS# XY2969000 IRRITANT	5g	41.70
		10g	51.90
		50g	205.50
A8,160-9	3-Amino-1,2,4-triazole, 95% [61-82-5] FW 84.08 mp 150-153° Beil. 26.137	25g	16.80
★	Merck Index 13,489 FT-IR 1(2).627A Safety 2,229A R&S 1(2).2367F RTECS# XZ3850000 CANCER SUSPECT AGENT MUTAGEN	100g	44.00
		1kg	188.20
A8,180-3	4-Amino-1,2,4-triazole, 99% [584-13-4] FW 84.08 mp 84-86° Beil. 26.16	5g	19.40
	FT-IR 1(2).627C R&S 1(2).2367H RTECS# XZ3850200 IRRITANT	25g	64.40
		100g	133.50
28,207-3	3-Amino-1,2,4-triazole-5-carboxylic acid hydrate, 98% [304655-78-5] FW 128.09 mp 182° Beil. 26.311 FT-NMR 1(3).91A Safety 2,229B R&S 1(2).2369B IRRITANT	25g	21.00
★	3-Amino-1,2,4-triazole-5-thiol, see 14,026-0. 3-Amino-5-mercapto-1,2,4-triazole page 91	100g	55.40
39,403-3	3-Amino-1-(2,4,6-trichlorophenyl)-2-pyrazolin-5-one, 97% [27241-31-2]	5g	14.80
★	FW 278.53 mp 218°(dec.) IRRITANT	25g	49.10
22,562-2	4-Amino-3,5,6-trichloropicolinic acid, tech. [1918-02-1] FW 241.46 mp 200°(dec.) Merck Index 13,7482 FT-IR 1(2).790B Safety 2,229C R&S 1(2).2543D RTECS# TJ7525000 CANCER SUSPECT AGENT	1g	21.50
40,307-5	4-Amino-2,3,5-trichloropyridine, 98% [28443-69-8] FW 197.45 mp 153-155° Beil. 22.433 IRRITANT	250mg	16.60
	2-Amino-1,1,3-tricyanopropene, see 10,741-7. 2-Amino-1-propene-1,1,3-tricarbonitrile page 106	1g	45.80
36,493-2	3-Amino-2,5,6-trifluorobenzoic acid, 98% [133622-65-8] H ₂ NC ₆ H(F) ₃ CO ₂ H FW 191.11 mp 175-180° FT-NMR 1(2).1156B R&S 1(2).1825A IRRITANT	1g	36.70
36,573-4	2-Amino-4-(trifluoromethyl)benzenethiol hydrochloride, 97% [4274-38-8] H ₂ NC ₆ H ₃ (CF ₃)SH·HCl FW 229.65 mp 196-198° FT-NMR 1(2).518C R&S 1(1).1417L IRRITANT	5g	34.90
		25g	115.90
24,892-4	7-Amino-4-(trifluoromethyl)coumarin, 99+% [53518-15-3] (coumarin 151) FW 229.16 mp 221-222° λ _{max} 207nm Safety 2,229D R&S 1(2).1947D IRRITANT	250mg	20.90
★	Suitable as laser dye	1g	57.80
19,696-7	2-Amino-5-trifluoromethyl-1,3,4-thiadiazole, 97% [10444-89-0] FW 169.13 mp 225-227° FT-NMR 1(3).118B FT-IR 1(2).651B Safety 2,230A R&S 1(2).2393H IRRITANT	1g	21.20
		5g	71.00
		25g	280.90
36,443-6	2-Amino-3,5,6-trifluoroterephthalonitrile, 98% [133622-66-9] H ₂ NC ₆ F ₃ (CN) ₂ FW 197.12 mp 138-140° FT-NMR 1(2).1543C R&S 1(2).2143G IRRITANT	1g	24.60
		5g	77.00
27,811-4	3-Amino-2,4,6-triiodobenzoic acid, 98% [3119-15-1] H ₂ NC ₆ H(I) ₃ CO ₂ H FW 514.83 mp 197-199° Beil. 14.414 FT-NMR 1(2).1156C Safety 2,230B R&S 1(2).1825B RTECS# DG3349000 IRRITANT	5g	28.40
		25g	94.50
44,436-7	5-Amino-2,4,6-triiodoisophthalic acid, 95% [35453-19-1] H ₂ NC ₆ I ₃ 1,3-(CO ₂ H) ₂ FW 558.84 mp 265-270° Beil. 14(4).1903 MOISTURE-SENSITIVE IRRITANT	10g	13.50
	May contain up to 5% water	50g	44.20
11,818-4	5-Amino-1,3,3-trimethylcyclohexanemethylamine, 99+% mixture of cis and trans [2855-13-2] (isophorone diamine) H ₂ NC ₆ H ₇ (CH ₃) ₃ CH ₂ NH ₂ FW 170.30 mp 10° bp 247° n _D ²⁰ 1.4880 d ₄ ²⁰ 0.922 Fp >230°F (110°C) [α] _D ²⁰ 0° (neat) CORROSIVE	250mL	21.90
★		1L	59.80
55,325-5	(S)-(-)-2-Amino-1,1,2-triphenylethanol, 97% [129704-13-8] C ₆ H ₅ CH(NH ₂)C(OH)(C ₆ H ₅) ₂ FW 289.38 mp 128-132° [α] _D ²⁰ -236° (c=1, CHCl ₃) HYGROSCOPIC	1g	103.60
★	A Product of Onyx Scientific, U.K.	5g	344.70

